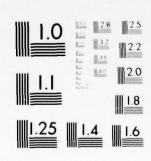


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∞ TRAP-ML-A TWO-DIMENSIONAL THERMAL RESPONSE CODE TAILORED FOR THE DEFENSE NUCLEAR AGENCY TRI-SERVICE THERMAL RADIATION TEST FACILITY Kaman AviDyne

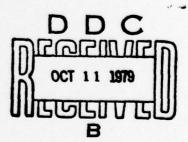
Kaman AviDyne 83 Second Avenue Burlington, Massachusetts 01803

30 November 1978

Final Report for Period November 1977-November 1978

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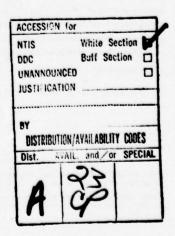
20. ABSTRACT (Continued)

The program is capable of analyzing multi-material, multi-layer samples exposed to either quartz-lamp radiation or a user-designated thermal source (as a function of both wavelength and time), with optional convective (air stream) cooling. Specific heats and thermal conductivities of the materials can be specified, in general, as functions of both temperature and radiation wavelength.

For output, the code computes the heat flow and temperature distribution throughout the finite model, and will optionally plot the temperature-time histories at layer boundaries and at both surfaces of the specimen.

PREFACE

The work was performed for the Defense Nuclear Agency (DNA) under Contract Number DNA001-78-C-0057 by Kaman Avidyne, Burlington, Mass., a division of Kaman Sciences Corporation. Captain J. M. Rafferty of DNA, and Major J. Hurst and Mr. G. Schmitt of the Air Force Materials Laboratory (AFML) at Wright-Patterson Air Force Base, Dayton, Ohio, were technical monitors. Mr. William Lee was project leader of the work performed in the Structural Mechanics Section of Kaman AviDyne headed by Mr. Emanuel S. Criscione, with technical contributions by Dr. J. R. Ruetenik and Dr. N. P. Hobbs.



Conversion factors for U.S. customary to metric (SI) units of measurement.

To Convert From	To	Multiply By
angstrom	meters (m)	1.000 000 X E -10
atmosphere (normal)	kilo pascal (kPa)	1.013 25 X E +2
bar	kilo pascal (kPa)	1.000 000 X E +2
barn	meter ² (m ²)	1 000 000 X E -28
British thermal unit (thermochemical)	joule (J)	1.054 350 X E +3
calorie (thermochemical)	joule (J)	4.184 000
cal (thermochemical)/cm ²	mega joule/m ² (MJ/m ²)	4.184 000 X E -2
curie	giga becquerel (GBq)*	3.700 000 X E +1
degree (angle)	radian (rad)	1.745 329 X E -2
degree Fahrenheit	degree kelvin (K)	t_ = (t* f + 459.67)/1.
electron volt	joule (J)	1.602 19 X E -19
erg	joule (J)	1.000 000 X E -7
erg/second	watt (W)	1.000 000 X E -7
foot	meter (m)	3.048 000 X E -1
foot-pound-force	toule (J)	1.355 818
gallon (U.S. liquid)	meter ³ (m ³)	3.785 412 X E -3
inch	meter (m)	2.540 000 X E -2
jerk	joule (J)	1.000 000 X E +9
joule/kilogram (J/kg) (radiation dose absorbed)	Grav (Gy)**	1,000 000
kilotons	terajoules	4.183
kip (1000 1bf)	newton (N)	4.448 222 X E +3
kip/inch² (ksi)	kilo pascal (kPa)	6.894 757 X E +3
ktap	newton-second/m ² (N-s/m ²)	1.000 000 X E +2
micron	meter (m)	1.000 000 X E -6
mil	meter (m)	2.540 000 X E -5
mile (international)	meter (m)	1.609 344 X E +3
ounce	kilogram (kg)	2.834 952 X E -2
pound-force (1bf avoirdupois)	newton (N)	4.448 222
pound-force inch	newton-meter (N·m)	1.129 848 X E -1
pound-force/inch	newton/meter (N/m)	1.751 268 X E +2
pound-force/foot ²	kilo pascal (kPa)	4.788 026 X E -2
pound-force/inch ² (psi)	kilo pascal (kPa)	6.894 757
pound-mass (1bm avoirdupois)	kilogram (kg)	4.535 924 X E -1
pound-mass-foot (moment of inertia)	kilogram-meter ²	
pound-mass/foot 3	(kg·m²) ktlogram/meter³ (kg/m³)	4.214 011 X E -2 1.601 846 X E +1
rad (radiation dose absorbed)	Gray (Gy)**	1.000 000 X E -2
roentgen	coulomb/kilogram (C/kg)	2.579 760 X E -4
shake	second (s)	1.000 000 X E -8
slug	kilogram (kg)	1.459 390 X E +1
torr (mm Hg, 0° C)	kilo pascal (kPa)	1.333 22 X E -1

^{*}The becquerel (Bq) is the SI unit of radioactivity; 1 Bq = 1 event/s. *The Gray (Gy) is the SI unit of absorbed radiation.

A more complete listing of conversions may be found in "Metric Practice Guide E 380-74," American Society for Testing and Materials.

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SECTION 1 INTRODUCTION

TRAP-ML is the result of a dual effort to (1) simplify the TRAP computer code (Reference 1) for studying radiation effects on aircraft components and (2) provide an analytical predictive and correlative tool for use in the Defense Nuclear Agency (DNA) Tri-Service Thermal Radiation Test Facility located at the Air Force Materials Laboratory, WPAFB, Dayton, Ohio. The appendage "ML" on the code name indicates the special applicability of the code to the Materials Laboratory.

The DNA thermal test facility, described in detail in References 2, 3 and 4, includes a test setup consisting of a quartz lamp bank for thermal radiation and a wind tunnel providing convective cooling on material samples approximately four inches square. The samples are mounted flush with the inside surface of the wind tunnel, facing a window in the opposite tunnel wall. The sample is irradiated through this window, simulating the environment of an aircraft exposed to nuclear radiation while in flight.

The TRAP code (Thermal Response of Aircraft to Nuclear Radiation) includes the material modelling and heat transfer capabilities required for the test facility application. The solution method is based on a finite representation of the cross section by thermally thin "elements." One or more elements make up a "layer," characterized by uniform material properties and thickness, and several layers are possible. The layers make up a "segment"; several segments are allowed, introducing the possibility of two-dimensional heat flow. The heat transfer mec sisms in TRAP include radiation, convection, and conduction.

The problem with the original TRAP is that it is too general for the test facility application. For example, it also includes a nuclear source generator, atmospheric transmission and ground reflection factors, and a vulnerability assessment including an iteration on slant

range. Also, it is designed to analyze large complex aircraft structures not laboratory samples. Clearly, these capabilities are not needed
for the test facility. Nor are the stress-strain calculations needed,
as the mechanical boundary conditions are not prescribed in the laboratory
setup.

Certain other capabilities were required in TRAP-ML, however, which are not in the original TRAP. The quartz lamp radiation source was modelled as a function of wavelength and time and included in the program. Provision was also made for two other source options, including a point-by-point option.

While the test facility setup can be treated as a one-dimensional thermal problem, the capability to analyze two-dimensional heat flow remains. For example, the airstream does tend to "heat up" downstream, introducing a nonuniform axial temperature field. So, rather than eliminate the 2-D capability, it was retained for possible future applications.

Also, while convection and reradiation are still important, these heat transfer processes were modified slightly to better simulate the test facility. Input and output were also modified to enable the user to operate the code with less than a $60,000_8$ core requirement and receive temperature-versus-time plots at each surface and layer boundary of the material sample.

TRAP-ML, then, allows the user to simulate a multi-layer, multi-material sample subjected to a general radiation source (a 6000 watt tungsten lamp in particular), complete with two-dimensional heat flow, convective cooling (if needed), and reradiation, with both the source and material absorptivity specified, in general, as functions of wavelength.

With this capability, the pretest calculations can be performed which, in many cases, may reduce the number of test exposures required. And where material properties are uncertain, post-test correlation with tests involving these materials may lead to evaluations of the unknown properties.

Section 2 describes in more detail the heat transfer processes in TRAP-ML, and Section 3 documents the computer program. Section 4 includes a sample problem which is compared with a test exposure of a .032 inch aluminum sample coated with 3 mils of white polyurethane. This limited correlation indicates excellent agreement. Finally, Table 1 includes a conversion table of units (English to Metric), since the TRAP computer code (and hence, TRAP-ML) was originally formulated in the English system of units, including input and output. Program listings make up Appendices A and B.

Table 1. Conversion of units (English to metric).

Value Expressed in English Units	Multiplied By	Yields Value Expressed In Metric Units
Inch (in)	2.54	Centimeters (cm)
British Thermal Unit (BTU)	252.0	Calories (cal)
BTU/in ²	39.03	Cal/cm ²
Temperature, °R	0.5555	Temperature, °K
BTU/in ² /°R	70.25	Cal/cm ² /°K
BTU/pound mass (BTU/1b-m)	0.5555	Cal/gram
BTU/1b-m/°R	1.0	Cal/gram/°K

SECTION 2 HEAT TRANSFER MECHANISMS

2-1 RADIATION SOURCE

The primary radiation source included in the program is the quartz lamp bank consisting of 6000 watt tungsten filament bulbs. The spectral distribution and a general time history are included in a non-dimensional manner, such that the user need only specify the peak flux level to be achieved and the time of exposure before the lamps are shut off. Provision is also made for other source characteristics, or even very simple sources such as a step-on and step-off time history.

The nominal spectral distribution is based on spectrometer measurements and the point-by-point model is indicated in Figure 1. While there is some absorption of the radiation by the atmosphere, it is small since the lamps are normally located within about 2 inches of the taermal sample, and is, therefore, neglected. The distribution is considered invariant with time, although it is thought the peak intensity may shift to a higher wavelength when cool. While this is ignored in the nominal source option, the capability exists of specifying the source as a function of both wavelength and time.

The program numerically integrates the distribution between 0.2 and 5.5 microns to determine the total flux, then normalizes the entire distribution to that value as a function of 14 discrete wavelengths. If wavelengths other than the nominal 14 are specified on input, the program linearly interpolates between the inputted values to determine values for the 14 basic wavelengths. These fourteen wavelengths and their associated band widths are indicated in Table 2. When material absorptivity is specified as a function of wavelength, these 14 bands are used to numerically integrate the total absorbed radiation between 0.2 and 5.5 microns. Since all the incident radiation

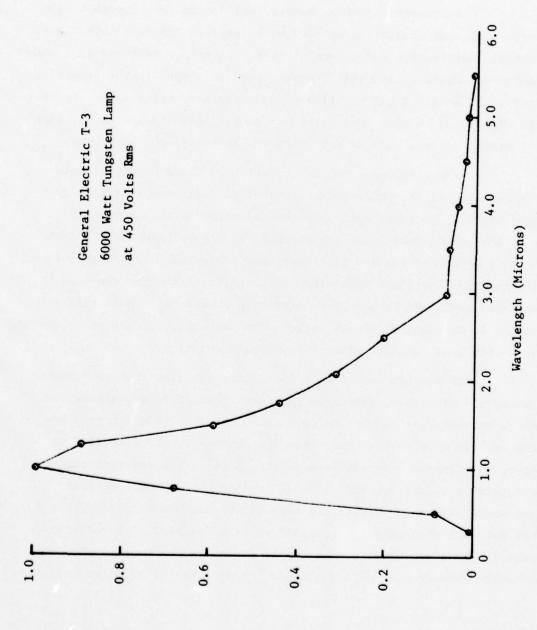


Figure 1. Spectral intensity model for nominal radiation source.

Relative Intensity, \overline{I}

Table 2. Wavelengths used to determine absorbed flux.

Wavelength Number	Nominal Wavelength (Microns)	Wavelength Band (Microns)
1	0.288	0.2 - 0.375
2	0.5	0.375 - 0.625
3	0.75	0.625 - 0.875
4	1.0	0.875 - 1.125
5	1.25	1.125 - 1.375
6	1.5	1.375 - 1.625
7	1.75	1.625 - 1.875
8	2.0625	1.875 - 2.25
9	2.5	2.25 - 2.75
10	3.0	2.75 - 3.25
11	3.5	3.25 - 3.75
12	4.0	3.75 - 4.25
13	4.5	4.25 - 4.75
14	5.0	4.75 - 5.25
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is either absorbed or reflected back into space, the incident absorbed relative intensity can be given by

$$\frac{1}{\overline{1}} \text{ abs} = \frac{0.2 \int_{0.2}^{5.5} \alpha(\lambda) \overline{1}(\lambda) d\lambda}{\int_{0.2}^{5.5} \overline{1}(\lambda) d\lambda}$$
(1)

where α represents the fractional absorptivity (0 $\leq \alpha \leq$ 1) and λ is the wavelength.

Figure 2 indicates the nominal time-wise radiation history, where the flux has been normalized to an arbitrary value. The time history is divided into two domains:

- The "rise" domain includes a very sharp rise in flux and then reaches a nearly steady-state value for large times. After 3.87 seconds the program assumes a constant flux.
- 2) The decay phase commences when the lamps are turned off and has a rapid decrease in flux until it finally is assumed to be zero at 2.85 seconds after the exposure has ended.

Regardless of when the lamps are turned off, the entire curve is normalized to a value of peak flux (q_{max}) which is specified on input. And the decay phase begins exactly with the time specified on input. The incident absorbed flux can then be given by

$$q^{abs}(t) = q_{max} \cdot \overline{1}^{abs} \cdot \overline{F}(t)$$
 (2)

where $\overline{\mathbf{F}}$ is the relative flux defined in Figure 2 and t represents time.

The user can also specify a simple step-on and step-off time history or a general point-by-point description. Linear interpolation is used between tabulated values. Even these distributions are normalized to the largest tabulated value so that the intensity of the radiation can be varied with a single input parameter. The entire specimen is assumed to be irradiated uniformly.

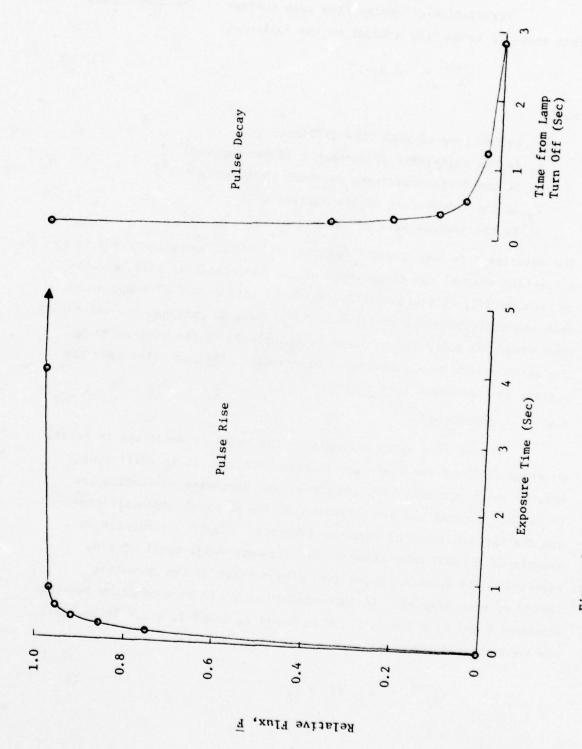


Figure 2. Time-wise exposure model of nominal radiation source.

2-2 RERADIATION

Reradiation of energy from each surface of the sample back into space is taken into account by the following:

$$q_i^{rad} = -A_i \varepsilon_i \sigma T_i^4$$
 (3)

where

q, is the rate of heat flow (BTU/sec)

 ϵ_i is the emissivity of surface i (dimensionless)

o is the Stefan-Boltzmann constant (BTU/sec-in²-°R⁴)

T, is the temperature of the surface (°R)

A_i is the surface area of element i (in²)

The emissivity is considered "hemispherical total" emissivity and is a function only of the temperature of the material. As will be noted in Section 3-1, if the absorptivity (α) is independent of temperature, then the emissivity can be taken as $1-\alpha$. But, in general, the emissivity must either be specified on input or calculated by the program using the absorptivity as a function of wavelength. The heat flow from the surface is considered lost forever.

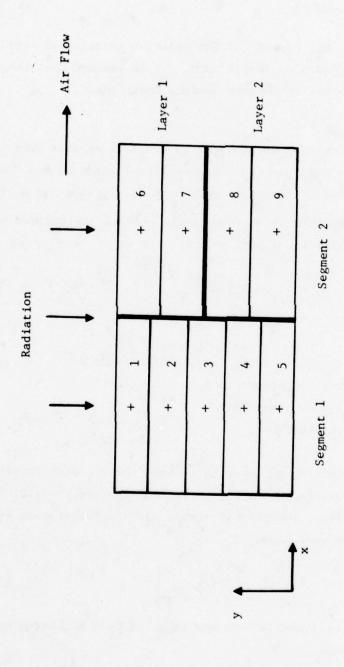
2-3 CONDUCTION

Conduction among elements in the model is described in detail on pages 86-92 of the TRAP report (Reference 1). It is sufficient here to point out that both lengthwise and depthwise conduction are taken into account and are functions of the material conductivities and the exposed areas of adjacent elements. Figure 3 indicates an example of a model consisting of two segments and a total of nine elements. The element "nodes" are always taken at the geometric center of each element. In this case there would be conduction between elements 1 and 6, 2 and 1, 2 and 6, 2 and 7, 2 and 3, etc. In general, the conduction between two elements is given by

$$q_{ij}^{cond} = \overline{K}_{ij} (T_i - T_j)$$
 (4)

+ Locates center of element.

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General two-dimensional cross-sectional model indicating elements, layers and segments. Figure 3.

where

K_{ij} is a conductivity factor between elements i and j, BTU/sec-°R
T is temperature, °R

As indicated earlier, \overline{K}_{ij} depends on the relative geometrical configuration of elements i and j, and also temperature. It is assumed that there are no conductive heat losses to the surrounding environment.

2-4 CONVECTION

Convective cooling is provided by means of exhaust fans at the outlet end of the wind tunnel. The model is located in the "throat" of the tunnel, a region of constant cross section, 1" wide and 4 1/2" high.

The boundary-layer heat flux to the surface is computed using the Newton heat-transfer equation

$$q^{conv} = h (T_{v} - T_{w}) A$$
 (5)

where

h is the heat transfer coefficient, BTU/in²/sec/°R

Tr is the recovery temperature, °R

Tw is the wall temperature, °R

A is the surface area, in²

Before discussing h, it is noted that the recovery temperature is computed with reference to the stagnation temperature, T_0 (which in this case is the ambient, room temperature), and the temperature at the outer edge of the body layer, T_{δ} ,

$$T_r = T_{\delta} + r \left(T_{\Omega} - T_{\delta}\right) \tag{6}$$

The recovery factor is dimensionless and taken to be 0.88, based on the discussion in Reference 1.

The free-stream temperature, \mathbf{T}_{δ} , can also be computed from \mathbf{T}_{o} , using perfect gas relations, where \mathbf{a}_{t} represents the total speed of sound in the room.

$$a_t = \sqrt{1.4 (1716) T_0} = \sqrt{2402.4 T_0}$$
 (7)

$$T_{\delta} = T_{o} \left\{ 1 - 0.2 \left(V/a_{t} \right)^{2} \right\}$$
 (8)

V is the free-stream velocity in ft/sec. The free-stream pressure, which will be needed later, can also be computed assuming a perfect gas,

$$P_{\delta} = P_{o} \left\{ 1 - 0.2 \left(V/a_{t} \right)^{2} \right\}^{3.5}$$
 (9)

where P is the ambient (room) pressure.

This, then, leaves the heat transfer coefficient, h, to be determined. This parameter is a function of many factors, but depends primarily on the type of flow in the region of interest, namely laminar vs. turbulent. This, in turn, depends on the tunnel geometry, roughness of the surface, any leaks in the tunnel, etc. Ideally, one probably would predict laminar flow in the test facility setup, but the discussion in Reference 4 indicates the flow is actually fully developed turbulent. Due to the factors mentioned above, this is plausible and accepted.

Based on turbulent flow, then, the Colburn equation (Reference 5) is used to take into account dependence of h on wall temperature and air-stream velocity (two velocities are possible, depending on whether one or two blowers is used). This requires either knowledge of the effective origin of the turbulent flow, or some empirical information regarding the value of h for a particular case. Reference 4 provided the latter, where an average value of $1.67 \times 10^{-4} \, \mathrm{BTU/sec/in^2/°R}$ was determined for an air flow of 793 ft/sec. The effective origin (x_{eff}) was then back figured to be 1.15 inches.

The Colburn equation and other necessary relations are presented below. The user is referred to Reference 1 for a more complete discussion.

$$h = \frac{0.026 \cdot k \cdot R_{eff}^{.8}}{x_{eff}}$$
 (10)

where

$$k = \left[1.0 - 8.89 (10^{-6}) \text{ T}^2 + 3.58 (10^{-8}) \text{ T}^2\right]$$

$$\cdot \left[2 (10^{-8}) \text{ T}^{-1.5} / (\text{T}^2 + 201.6)\right]$$
(11)

$$R_{e_{eff}} = \frac{R_{e} \cdot V \cdot x_{eff}}{12}$$
 (12)

$$R_{\delta} = 3.7 (10^6) P_{\delta} (T' + 198.6)/(T')^{2.5}$$
 (13)

$$T' = 0.5 (T_W + T_{\delta}) + 0.22 (T_r - T_{\delta})$$
 (14)

The program also provides the user with the optional capability of specifying the heat transfer coefficient, h, which allows him more direct control over that parameter.

2-5 TEMPERATURE CHANGE

The computer program combines the results of all the heat transfer mechanisms at each instant of time for each element. The temperature change can then be calculated for the finite time interval according to

$$\Delta T = \frac{\Delta t}{\rho A c_p} q \tag{15}$$

where

AT is the temperature rise, °R

At is the time interval, sec

o is the element density, lbs/in3

c is the specific heat of the material, BTU/1b-°R
A is the cross-sectional area (volume) of the element, in²
q is the net heat flow into the element, BTU/sec.

SECTION 3 COMPUTER PROGRAM DESCRIPTION

3-1 PROGRAM OPERATION

The TRAP-ML program was written in FORTRAN IV and developed on the Control Data Corporation (CDC) 6600 computer. The program requires card input, which is described in Section 3-2, and produces time history response output, the subject of Section 3-3. The program also has the capability to produce a plot file of temperatures versus time. This data file then serves as input to a second program, APLOT, which produces paper plots. This second program requires no card input of its own.

To execute and also generate plots, then, requires the execution of both TRAP-ML and APLOT back to back. The primary reason for separating the two codes was to enable the user to run with less than a $60,000_8$ cell core requirement. This makes it possible to operate the code from remote terminal under current WPAFB restrictions. A typical job control setup is shown in Table 3.

The routines for both programs are listed in Table 4. The basic plot routines are included in the APLOT package, rather than calling on system routines, so that the program is less dependent on one particular computer.

TRAP-ML requires approximately 53,000₈ to load and execute. Logical files TAPE1 and TAPE6 are used for input and output, respectively. TAPE5 is used internally to store the input data subsequent to reading the data into the computer. The plot information is written to file TAPE8 and is automatically rewound prior to and upon completion of job. More than one case can be run during a job execution, in which case the plot output for each case represents a separate file on TAPE8. Approximately 8 cp seconds are required to compile TRAP-ML using "OPT=1" and "R=2" options.

Table 3. Typical job control sequence.

Job Control Cards
JOB CARD MAP,PART ATTACH,A,TRAPML A. ATTACH,B,APLOT B. ROUTE,PLOT, EXIT. 7/8/9 input data for TRAPML

Table 4. TRAP-ML and APLOT routines.

TRAP-ML	APLOT
TRAPML	APLOT
BLOCK DATA	FIRST
CONVEC	KALINE
DTEMP	KALE
FLUX	KANUM
INT1Z	KASYM
RPLOT	KAVANS
RPRINT	KAXIS
SETUP	
SPECT	
TIN	
TSTEP	
XHEAT	

Program APLOT only requires the file TAPE8 for input, but does generate output on TAPE6. The program also ultilizes file TAPE40 for internal purposes. If on-line plots are requested, then the file PLOTS represents the resultant plot file, ready to be directed to a plotter.

APLOT requires approximately $50,000_8$ cells of memory to load and execute. Compilation requires only about four cp seconds.

Execution times will depend on the size of the model, the time interval required, and the duration of the response requested, but generally program APLOT will execute for only a small fraction of the time used by TRAP-ML. The example problem discussed in Section 3-4 required about 40 cp seconds for both programs.

3-2 PROGRAM INPUT

TRAP-ML looks for BATCH input on file INPUT. The input data are specified in groups, where each group begins on a separate card. More than one card may be required for a group, however. The variable type and format corresponding to each data group is given in parentheses in the input instruction and is always in fields of 12. For convenience, floating point numbers can be left justified in the field as long as the exponent is right justified. Also, zero values can be replaced by a blank field. Columns 73 through 80 are not used for data and can be used for card identification or other purposes.

All input parameters, where appropriate, should be compared with the maximum dimensions provided for in the program, as delineated in Table 5. This is very important since the program does not attempt to check the input for all such violations. In Table 5, the input parameters are delineated with an asterisk (*) preceding the variable name.

The specific instructions for input are contained in Table 6.

The remainder of this section will attempt to amplify on those instructions and provide insight into the applicability and limitations of the program.

Table 5. Description of variables which determine program dimensions.

VARIABLE	LIMIT	DESCRIPTION
*NAT	10	Number of temperatures at which absorptivity is given.
NCONDC	122	Number of conduction connections between elements.
*NCPT	10	Number of temperatures at which specific heat is given.
*NCT	10	Number of temperatures at which conductivity is given.
NEL#	75	Total number of elements.
NFALL	. 8	Number of points at which the relative flux is specified versus time in the decay phase.
*NFLUX	15	Number of points describing thermal flux time history.
NLAY	5	Number of layers in segment.
NMAT	9	Number of different materials.
NRISE	8	Number of points at which the relative flux is specified versus time in the "rise" phase.
*NSEG	3	Number of segments.
*NTSPEC	3	Number of times at which spectral distribution of radiation source is specified.
NWL	14	Number of wavelength bands used to describe radiation source in calculations.
*NWLA	10	Number of wavelengths at which absorptivity is specified for new material.
*NWLI	50	Number of wavelengths used to describe spectral distribution.

NOTES:

- * Variable is specified on input.# Dimension also limited in subroutine SETUP by a fixed point number before statement number 23.

Table 6. TRAP-ML input.

Group 1: (40A2) (ID(1), I=1, 40)

Title card - free field. (ID)

Group 2: (2112) INOUT, IPLOT

Program output-option code: (INOUT)

0, do not print input data

1, print input data

Plot code (IPLOT)

0, do not generate plots of temperature vs. time

1, do generate plots of temperature vs. time

Group 3: (I12) NSEG

Number of segments. (1 \leq NSEG \leq 3)

Group 4: (2F12.1) XSEG(IS), YSEG(IS)

x-coordinate of end of segment, in (XSEG) y-coordinate of end of segment, in (YSEG)

Note - See Figure 3 for definition of coordinates.

Repeat Group 4 for IS=1, NSEG+1

Groups 5-7 provide data for each segment, IS, and are to be repeated for IS=1, NSEG

Group 5: (I12) NLAY(IS)

Number of layers in segment $(1 \le NLAY(IS) \le 5)$.

Groups 6-7 provide data for each layer, IL, and are to be repeated for IL=1, NLAY(IS)

Group 6: (F12.1) TLAY (IS, IL)

Thickness of layer, in. [TLAY(IS,IL)]

Group 7: (2112) KMAT (IS, IL), NELL(IS, IL)

Code defining material of layer: (KMAT (IS, IL))

1, Aluminum, 2024-T3

2, Aluminum, 7075-T6

3, Magnesium, AZ31B-H24

4, Titanium, Ti-8Mn

Table 6. (Continued)

If materials other than those listed above are used, they should be assigned numbers which are consecutive beginning with 5. Material properties for the added materials are called for in Groups 8-21

Number of elements in layer. (NELL(IS, IL))

NM, and should be repeated until all added materials have been completed. If there are no added materials, skip to Group 23.

Group 8: (112) MCOR(NM)

Number of material to which new material partially corresponds. This number may be any number lower than the number of the new material for which data are now being supplied. Use of a corresponding material allows the analyst to skip reading in material properties which are the same as those of the corresponding material. If this number is read in as zero, all material properties for the new material must be supplied, starting with Group 10 and ending with Group 22. (MCOR(NM))

If MCOR = 0, skip to Group 10.

Group 9: (I12) KCH

Code indicating material property to be changed:

- 1, absorptivity
- 2, conductivity
- 4, specific heat
- 5, melting temperature
- 6, heat of fusion
- 11, density

Under this mode of input, a change card is read in for each property to be changed, followed by data for the specified material property. Group 22 terminates this sequence. (KCH)

Groups 10-13 provide absorptivity data for the new material, if required.

(MCOR = 0 or KCH = 1)

Group 10: (I12) NAT(NM)

Number of temperatures at which absorptivity is given NAT(NM))

Table 6. (Continued)

Groups 11-13 provide data for one temperature, NT, and are to be repeated for NT = 1, NAT(NM).

Group 11: (2F12.1) TALF(NT, NM), EMI(NT, NM)

Temperature at which absorptivity is given, °R (TALF(NT,NM))

Emissivity at specified temperature, dimensionless (EMI(NT,NM)) (Omit if integration of absorptivity by program is desired in order to obtain emissivity)

Group 12: (I12) NWLA

Number of wavelengths at which absorptivity is specified (NWLA)

Group 13: (2F12.1) WL(NW,NT), ALF(NW,NT)

Wavelength at which absorptivity is given, microns* (WL(NW,NT))

Absorptivity, dimensionless (ALF(NW,NT))

Repeat Group 13 for NW = 1, NWLA.

Groups 14-16 provide conductivity data for the new material, if required (MCOR = 0 or KCH = 2)

Group 14: (I12) KLD(NM)

Code of conductivities: (KLD(KM))

- 1, lengthwise and depthwise conductivities are the same
- 2, lengthwise and depthwise conductivities are different

Group 15: (I12) NCT(NM)

Number of temperatures at which conductivity is given (NCT(NM))

Group 16: (3F12.1) TCOND(NT,NM), CONDL(NT,NM), CONDD(NT,NM)

Temperature at which conductivity is given, °R (TCOND(NT,NM)

Lengthwise conductivity of material, BTU/in-sec-°R
(CONDL(NT,NM))

^{*}If emissivity has been specified in Group 11, only wavelengths from 0.2 to 5.5 microns must be covered. If emissivity is to be found by integration of absorptivity, wavelengths from 0.2 to about 250 microns must be covered.

Table 6. (Continued)

Depthwise conductivity of material (Omit if KLD(NM) = 1), BTU/in-sec-°R (CONDD(NT,NM))

Repeat Group 16 for NT = 1, NCT(NM)

Groups 17-18 provide specific heat data for the new material, if required (MCOR = 0 or KCH = 4)

Group 17: (I12) NCPT(NM)

Number of temperatures at which specific heat is given (NCPT(NM))

Group 18: (2F12.1) TCP(NT, NM), CPM(NT, NM)

Temperature at which specific heat is given, °R (TCP(NT,NM))

Specific heat of material, BTU/1b-°R(CPM(NT,NM))

Repeat Group 18 for NT = 1, NCPT(NM)

Group 19 provides melting temperature data for the new material if required (MCOR = 0 or KCH = 5)

Group 19: (F12.1)

Melting temperature, °R (TMELT(NM))

If the melting temperature is changed, usually all of the material properties will have to be changed. The reason for this is as follows: the me ting temperature for each material for which data are proceeded in the program is the same as the maximum temperature for the material property data. Suppose that the melting temperature is increased. Unless the other me erial properties are also changed, the melting temperature will be above the maximum temperature at which material properties are available. The program will then produce a error stop.

Group 20 provides heat of fusion data for the w material, if required (MCOR = 0 or KCH = 6)

Group 20: (F12.1) HOF (NM)

Heat of fusion of material, BTU b (HOF(NM))

Table 6. (Continued)

Group 21 provides density data for the new material, if required (MCOR = 0 or KCH = 11)

Group 21: (F12.1) RHOM(NM)

Density of material, lbs/in (RHOM(NM))

Group 22: Blank card. Is required for each new material NM, either at the completion of Groups 10-21 for MCOR(NM) = 0, or at the completion of the sets of data beginning with Group 9 for NCOR(NM)70.

Group 23: (4F12.1) VEL, PO, TAM, HCONO

Velocity of free-stream flow, ft/sec (VEL)

Ambient (room) pressure, psi (PO)

Ambient (room) temperature, degrees F (TAM)

Convective heat transfer coefficient, BTU/in²-sec-°R (HCONO)

Note - Program computes coefficients for each segment if HCONO=0.0.

Group 24: (112) NOPT

Code designating thermal flux time history (NOPT)

1, 6000 watt tungsten source (Figure 2)

2, Step load.

3, Arbitrary point-by-point description.

Group 25: (F12.1) QDMAX

Maximum flux achieved, BTU/in²-sec (QDMAX)

If NOPT=3, skip Group 26

Group 26: (F12.1) TCUT

Time at which thermal source is cut off, sec. (TCUT)

If NOPT<3, skip Groups 27-28.

Group 27: (112) NFLUX

Number of points describing thermal flux time history. First point must include time=0. (2 < NFLUX < 15)

Table 6. (Continued)

Group 28: (2F12.1) TFLUX(I), FLUXT(I)

Time, sec (TFLUX(I))

Relative flux corresponding to TFLUX, BTU/in²/sec. Magnitude of flux is adjusted to QDMAX. (FLUXT(I))

Repeat Group 28 for I = 1, NFLUX

Group 29: (2112) NWLI, NTSPEC

Code designating spectral distribution of flux (NWLI)

0, 6000 watt tungsten characteristics for all times (Figure 1).

>0, number of wavelengths used to describe distribution (2<NWLI<50).

Number of times at which spectral distribution is specified. Assumed to be 1 if NWLI=0. ($1 \le NTSPEC \le 3$).

If NWLI=0 or NTSPEC=1, skip Group 30.

Group 30: (6F12.1) (TSPEC(I), I=1, NTSPEC)

Times at which spectral distribution is specified. First time must be 0.0. (TPSEC)

If NWLI=0, skip Group 31.

Group 31: (2F12.1) (WAVE(J,I), QR(J,I))

Wavelength at which flux is specified. First wavelength must be 0.2 microns; the last must be 5.5. (WAVE)

Relative flux corresponding to WAVE, BTU/in²/sec. Only relative magnitudes are important. (QR)

Repeat Group 31 for J=1, NWLI, then repeat that sequence for I=1, NTSPEC.

Group 32: (3F12.1) DELTIM, TSTOP, PRINT

Integration time interval (if read in as 0.0, program will determine time interval required for stability), sec (DELTIM)

Time at which computations are to stop, sec (TSTOP)

Number of time intervals between printouts (0.0 will give no printout), dimensionless (PRINT)

Table 6. (Concluded)

To execute another problem, repeat Groups 1-32.

Group 33: (A3) IEND

To terminate job, place "END" in columns 1-3. (IEND)

TRAP-ML has the capability to assemble a two-dimensional heat transfer model of a particular material specimen. This implies that the specimen and the thermal loads are uniform in one coordinate direction - defined here to be the z direction. The x-coordinate direction is defined as in the direction of the air flow, and y is defined to be from the unexposed surface through the cross section toward the exposed surface (see Figure 3).

The model consists of segments, layers and elements, where elements are the basic "building blocks" and must be thermally "thin"; i.e., the temperature gradient is accurate only to the extent that enough elements are included in the model. Geometrically, the elements are horizontal slices of the cross section of constant thickness. Layers consist of a set of elements, all of which exhibit identical material properties. The model can contain up to five layers in the cross section per segment. The segment is defined as a depthwise set of layers representing a rectangular portion of the cross section. Elements, layers, and segments all have constant thickness. If only one segment is included in the model, the program is, by definition, reduced to a one-dimensional analysis. It is expected that only one segment will be used to model the test specimens.

In the general model indicated in Figure 3, segment 1 has only one layer and five elements. Segment 2 consists of two layers, each having two elements.

Now for the input instructions in Table 6. Group 1 represents a free-field case descriptor which appears as a title in the output. The first 40 columns will also appear on plots if requested. It should be pointed out that data groups 1 through 32 are repeated for as many cases as the user desires.

Group 2 defines the output options. By selecting INOUT=1, the output is reduced somewhat and the user still gets a raw data listing at the beginning of the output. The plots which are generated if IPLOT=1 give temperature at both surfaces and at each layer boundary versus time.

Group 4 defines the location and size of each segment by specifying the lower, left-hand and right-hand coordinates of each segment as defined in Figure 3. The coordinate origin can be located anywhere - one might just as well start with (0., 0.) for the first segment as not. Groups 5 and 6 define the layers.

Group 7 introduces the material properties. If materials or material properties other than those associated with the four materials indicated are desired, the user can accomplish this by calling for one or more additional materials. The first new material is number 5, the second number 6, etc. The properties corresponding to the four built-in metal alloys are indicated in Table 7 and Figures 4 and 5. In this regard, it will be noted that the selection of these four materials was arbitrary and the user is referred to Reference 1 for a complete discussion of property specification. It should be noted, for example, that the absorptivities specified for materials 1-4 are all 0.5 for an unpainted condition. If the user wants to specify a different absorptivity, he would need to call for a new material and proceed with the input.

Also in Group 7 the user must decide on the number of elements per layer. The program automically divides each layer into elements of equal thickness, depending on the number (NELL) specified. In general, thicker layers with lower conductivities will require more elements. One guideline, other than checking the resulting output to see that the temperature gradient is not too large, is to check the Δt calculation (see the discussion of Δt in Group 32). Here it makes some sense to end up with a Δt requirement more or less the same for all elements. At least two elements per layer are recommended so that a good estimate of temperature can be made at the boundary between layers and at the surface.

Group 32 requires the selection of a Δt to be used in the numerical integrations. This number must be small enough to give a stable solution. If Δt =0.0 is specified, the program will compute a Δt associated with each element and then select the smallest of these.

Table 7. Material properties in TRAP-ML.

Material	Aluminum 2024-T3	Aluminum 7075-T6	Magnesium AZ31B-H24	Titanium Ti-8Mn
Absorptivity	0.5	0.5	0.5	0.5
Emissivity	0.5	0.5	0.5	0.5
Heat of Fusion (BTU/1b)	170.	170.	158.	188.
Density (1b/in ³)	0.100	0.101	0.064	0.171
Melt Temp. (°R)	1680.	1680.	1662.	3495.

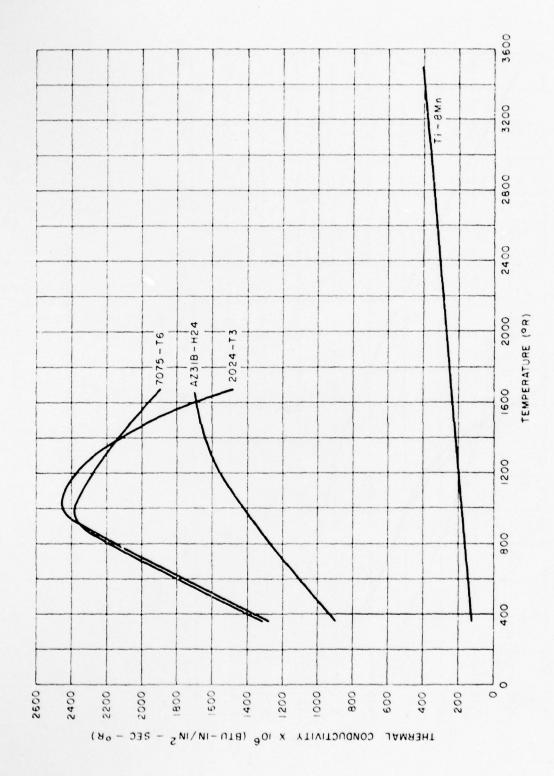


Figure 4. Thermal conductivity versus temperature.

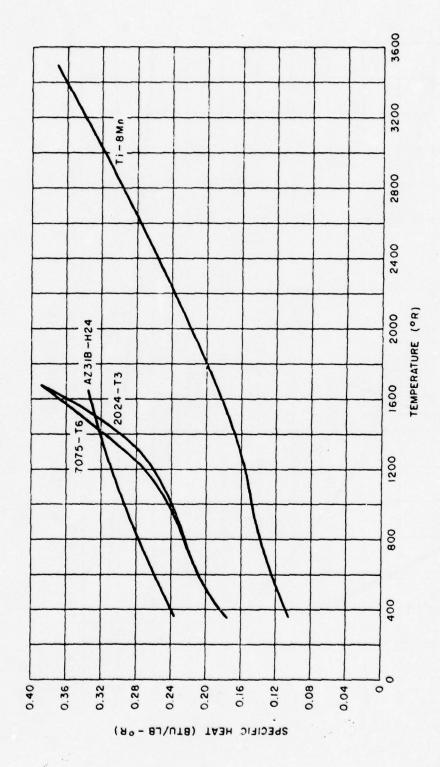


Figure 5. Specific heat versus temperature.

The user is referred to Reference 1 for a complete discussion of the formulation, but the resultant formula used is the following:

$$\Delta t_{i} = \frac{0.8 \, \rho_{i} \, A_{i} \, C_{p_{i}}}{\sum_{j} \overline{K}_{ij}}$$
(16)

where the summation on j extends over all elements which touch element i. These numbers are then printed out for the user's information.

Groups 8 through 22 permit the user to specify new material properties. Note that if the new material (called for in Group 7) does not have totally different properties from one of the four basic material options, then only those properties that change need be specified. For example, if one were to specify the absorptivities for 7075-T6 Al, then MCOR (Group 8) would be 2 and KCH (Group 9) would be 1. The absorptivities would be specified in Groups 10-13.

With regard to melt temperature (Group 19), the program treats the melting of an element by eliminating it from the solution. The next element then becomes exposed to the radiation source, with the implicit assumption that the melting always proceeds from the exposed surface toward the opposite surface, element by element. The output reflects this event by setting the temperature of a melted layer to 0.0 in the printed output, and a small vertical arrow indicates the event on the plotted output.

3-3 PROGRAM OUTPUT

TRAP-ML output consists of three logically distinct sections. The first output is a card image listing of the entire input deck, exactly as it is read in, except that the cards are numbered. The second section describes the input data (if INOUT=1) in terms of its meaning in the program.

The third section is the time-history response output. At each printout time the following values are printed: time, incident flux, incident fluence, and a table of (1) temperatures and (2) a

breakdown of heat flow factors for each element. Figure 6 represents typical time-history output for three consecutive printouts.

If plots are requested (IPLOT=1), the program also stores between 100 and 200 sets of temperatures (evenly spaced in time) for both surfaces of the model and for each layer division for segment 1, or for segment 3 if there are 3 segments in the model. Linear interpolation/extrapolation is used to estimate these temperatures based on the available values at the center of each element. Each curve plotted is identified by a unique symbol. These symbols are defined in Table 8. The final two lines of printed output indicate the number of data points per curve and the number of curves written to disk file.

Error messages are largely self-explanatory, although the linear interpolution routine (INT12) only indicates a code number indicating the source of a problem if a variable exceeds the limit of the table. Table 9 identifies the routine corresponding to the error code.

Program APLOT generates a list of the data plotted and indicates the completion of each graph as it is actually generated.

NET .72665-02 .74665-02 .74625-02 .80795-02 .23925-02	NE1 • 6754E - 02 • 1351E - 02 • 7492E - 02 • 7492E - 02 • 2264E + 00	NET 5781E-02 5115E-02 355E-02 2361E-02 5515E-01
. S E C C C C C C C C C C C C C C C C C C	. SE C C C C C C C C C C C C C C C C C C	5 E C COMOUCTION .1207E+00 .5115E-02 .3365E-02 .2361E-02 .5515E-01
B T U	9 T U CONVECTION 2909E+00	6 T U CONVECTION2191E+00
AHSORPTION READIATION .7908E+008184F-02	.819455E+U0 .151765E+U1 H E A T F L D M ABSORPTION REMADIATION .7936E+UU1036E-01	.1015946+00 .1562976+01 H E A T F L O M ABSORPTIUM RERADIATION .98346-0156946-02
18 4 4 4 4 4 4 4 4 4 4 4 4 4 4 4 4 4 4 4	1940496+01 UENCE, BTU/IN**2 = EMENT TEMPERATURE 1056755 04 2 .92015+03 4 .92015+03 5 .94745+03 7 .97045+03	** ** ** ** ** ** ** ** ** ** ** ** **
ELEVENT TEAPERATURE NUMBER DEGREES R - 9407E+03 - 9185E+03 - 8686E+03 - 6595E+03 - 6595E+03 - 6595E+03 - 6595E+03 - 6595E+03	77 72 mx = 40 x 40 x 4	rr m2 3 m2 = 0 w = n v v = r 1 1 2 2 2 2 2 2 2 2
SECUENT EL	INCIDENT INCIDENT SEGMENT NUMBER I	INCIDENT INCIDENT INCIDENT SEGMENT NUMBER 1

Figure 6. Time-history output for sample problem.

Table 8. Plot symbols used to identify plot curves.

Curve Number (front face first- back face last)	Plot Symbol
1 (front face)	
2	Δ
3	×
4	♦
5	+
6 ↓	0

Note - A vertical arrow (♠) indicates the time at which the element melted.

Table 9. Error codes from INT1Z.

Error Code	Subroutine which called INT1Z	Location Code #
1	SPECT	300-2
2	FLUX	300
3	SPECT	700 ⁻¹
4	SETUP	228-3
5	TSTEP	12
6	TSTEP	13
7	DTEMP	7
8	DTEMP	10
9	XHEAT	62
10	XHEAT	83
11	DTEMP	25
12	SETUP	117 ⁻¹
13	TSTEP	12-4
14	XHEAT	42

 $\# The \ location \ code$ is read as follows: a code $S^{\mbox{+}n}$ is interpreted as FORTRAN statement number S plus or minus n cards.

SECTION 4

SAMPLE PROBLEM AND COMPARISON WITH EXPERIMENT

The sample problem described in this section is designed to (a) provide the user with a check case for exercising the code, (b) provide an example of input data preparation, and (c) make a limited comparison with experimental results. The example case selected is based on an experiment performed at the DNA test facility and documented in Reference 4. A four inch sample of 2024-T3 aluminum substrate .032 inches thick was coated with .003 inch of white polyurethane and subjected to a 793 ft/sec wind tunnel flow and a two second pulse from the quartz lamp bank.

The input data for TRAP-ML is indicated in Figure 7. Although the data corresponds to the test data as closely as possible, the documentation was not totally complete. Figure 8 represents the plot generated by APLOT of three temperature-time histories: (1) the exposed surface, which obtained a peak temperature of $1019^{\circ}R$, (2) the division between coating and substrate, and (3) the unexposed surface (backface), which obtained a peak temperature of $900^{\circ}R$. The last two curves fall virtually one on top of the other due to the relatively high conductivity of the aluminum. Remember, the plot symbol codes are tabulated in Table 8. The printed output in the vicinity of where peak front-face temperatures were computed can be found in Figure 6.

Figure 9 compares the backface temperature with that presented in Figure 26 of Reference 4. Also shown are the results of the ASTHMA code from Reference 4. In general, TRAP-ML seems to predict the temperature response well - its only shortcoming is a failure in this case to predict as large a peak temperature. This minor deviance could easily be explained by any number of uncertainties in the data, e.g., the exact shape of the radiation pulse time history.

```
WHITE POLYURETHANE OVER .032 IN. 2024-T3 SAMPLE PROBLEM.
0.
            0.
           5
.003
           5
                       5
.032
           1
                       2
           0
           5
460.
            .756
5.5
            .242
            .242
           .758
4000.
.2
5.5
            .242
            .242
           2.54
0.
                     E-6
            2.54
                     E-6
            .41
0.
4000.
            .41
1200.
50.
793.
                      75.
           14.3
                                0.
.82
2.
           0
                        500.
0.
END
```

Figure 7. Sample problem input data.

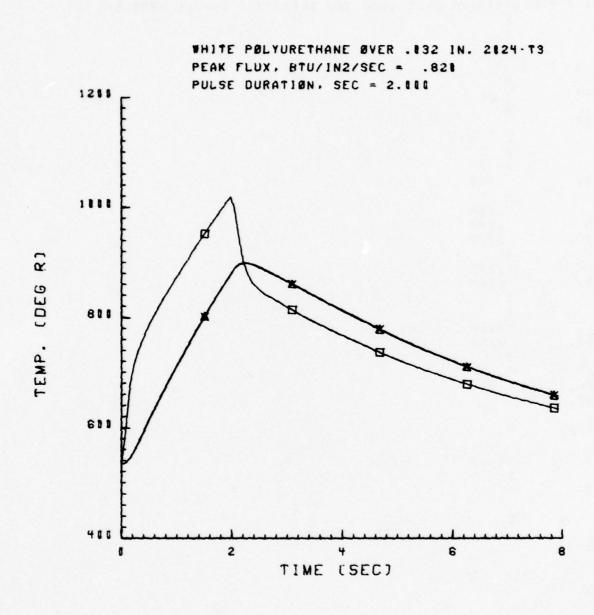


Figure 8. Computer-generated sample plot.

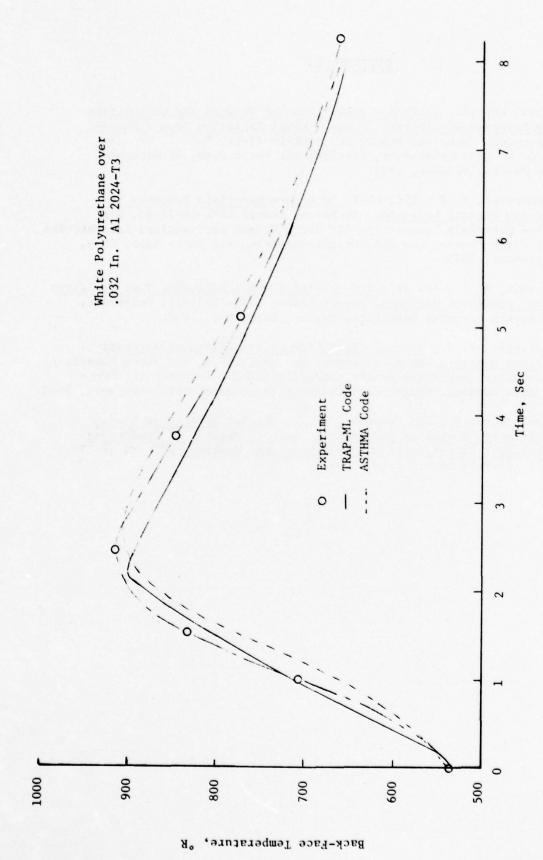


Figure 9. Comparison of TRAP-ML predictions to experiment.

REFERENCES

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- Servais, R. A., et. al., <u>Tri-Service Thermal Radiation Test Facility</u>: <u>Test Procedures Handbook</u>, Report Number UDRI-TR-77-28, University of Dayton Research Institute, Dayton, Ohio, May, 1977.
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 Coating Systems, Technical Report No. AFML-TR-78, Air Force Materials
 Laboratory, Air Force Wright Aeronautical Laboratories, Air Force
 Systems Command, Wright-Patterson Air Force Base, Ohio, February, 1978.
- 5. Blessing, A. H., and Hanawalt, A. J., Thermal Effects on Static Aeroelastic Stability and Control, Part I, Methods of Predicting Structural Temperatures Due to Aerodynamic Heating, WADC TR 58-378, Bell Aircraft Corp., June, 1959.

APPENDIX A
PROGRAM LISTING OF
TRAP-ML

```
PPOSPAM TRAPML (INPUT, OUTPUT, TAPE1=INPUT, TAPE6=OUTPUT, 1 TAPE5=129, TAPE8=514)
                           TRAP-ML IS A SPECIAL VERSION OF TRAP DEVELOPED FOR THE AIR MATERIALS LABORATORY, WPAFB, BY KAMAN AVIDANE, BURLINGTO VERSION 1.0, SEPT., 1978.
TEMPERATURE VS. TIME PLOT INFORMATION IS WRITTEN TO TAPES.
                                                                                                                             VERSION OF TRAP DEVELOPED FOR THE AIR FORCE
                     COMMON A (75).ALF (10,16).B(122).CKD(75).CKDM(9).CKL(75).CKLM(9).

2 ONDAG (122).CONDD(10,9).CONDI(122).CONDJ(122).CP(75).CPMAT(9).

3 IELTIM.DTIM(75).DTIME.DTMIN.EL(75).FLUXT(15).HCONO.ID(20).

4 IEL(3).IELAY(6,3).IEND.IMAT(75).IMELT(75).INOUT.IPLOT.ISEG(75).

5 ISTOP, JEL(3).KCH.KCONDI(122).KCONDJ(122).KCPI(122).KCPJ(122).

6 NFLUX.NLAY(3).NMAT.NOPT.NPRINT.NSEG.NTSPEG.NM.A.NWLI.PO.

7 PRINT.D(75).JAB(3).DCOND(75).QCONV(3).QDMAX.JFLU.QR(50.3).

8 QRAD.QRER(2.3).OPR(14).SEGL.T(75).TAM.TCUT.TEF(3).TEMP(75).

COMMON TFLUX(15).TFS.TIME.TLAY(3.5).TPC(75).TR,TSPEC(3).TSTOP.

1 VEL.WAVE(50.3).WAVEL(14).WL(10.10).WT(75).X(75).XCON(75).

2 XSFG(4).XTR.Y(75).YSEG(4).

COMMON /BLOCK/ ALFAT(10.14.9).CONDL(10.9).CPM(10.9).EMI(10.9).

1 HOF(9).KLD(9).NAT(9).NCPT(9).NCT(9).NMATT.NWL.PI.RHOM(9).

1 TALF(10.9).TOONO(10.9).TCP(10.9).TMELT(9).WAVEB(15)
                           CALL TIN
                       CALL SETUP
IF (ISTOP.EQ.1) GC TO 100
IF (DELTIM.EQ.0.0) CALL TSTEP
CALL XMEAT (0)
DO 10 IE=1,NEL
TEMP1 | TAM
IMELITIE) = 0
O(IE) = 0.0
TIME = 0.
NPRINT = 0
OFLJ = 0.
                           CALL XHEAT (2)
CALL DTEMP(0)
IF (IPLOT.GT.C) CALL RPLOT(0)
3
            70 IF (IPLOT.GT.0) CALL RPLOT(1)
IF (NPRINT.GT.0) GO TO 75
IF (PRINT.EQ.0.8) GO TO 77
CALL PPRINT
NPRINT=PRINT-0.5
GO TO 77
75 NPRINT=NPRINT-1
77 TIME=TIME+DELTIM
IF (TIME.GT.TSTOP) GO TO 90
CALL XHEAT (2)
CALL DIEMP(1)
OFLJ = QFLU + QRAD*DELTIM
GO TO 70
             90 IF (IPLOT.GT.0) CALL RPLOT(2)
        100 ENDFILE
REALIND 8
120 STOP
END
```

TRAPHL

```
BLOCK DATA
                               COMMON /BLOCK/ ALFAT(10.14.9), CONDL(10.9), CPM(10.9), EMI(10.9), 
4DE(9), KLD(9), NAT(9), NCPT(9), NCT(9), NMATT, NHL, PI, RHOM(9), 
TALE(10.9), TCOND(10.9), TCP(10.9), TMELT(9), HAVEB(15)
                                ABSORPTIVITY.
                          DATA NAT/4*2/.ALFAT/560*0.5/.TALF(1.1)/300.0/.
1TALF(2.1)/1680.0/.TALF(1.2)/300.0/.TALF(2.2)/1680.0/.
2TALF(1.3)/300.0/.TALF(2.3)/1662.0/.TALF(1.4)/300.0/.
3TALF(2.4)/3495.0/
                                SPECIFIC HEAT. STUZER-DEG R.
                               DATA NCPT/9.9.3.9/
                       DATA NCPT/9,9,8,9/

MATERIAL 1, 2024-T3.
DATA CPM/0.175.0.191.0.214.0.229.0.244.

10.259.0.287.0.313.0.387.0.387.
MATERIAL 2, 7075-T6.
20.174.0.190.0.213.0.230.0.250.
30.241.0.307.0.332.0.389.0.389.
MATERIAL 3, AZ31R.
40.237.0.246.0.264.0.281.0.298.
50.312.0.324.0.336.0.336.0.336.
MATERIAL 4, TI-8MN.
60.104.1275.0.146.0.153.0.158.
70.154.0.204.0.251.0.370.0.370/
MATERIAL 1, 2024-T3.0.660.0.860.0.1060.0.
1250.0.1360.0.1460.0.1680.0.1680.0.
1250.0.1360.0.1460.0.1680.0.1680.0.
MATERIAL 2, 7075-T6.
2360.00.460.00.660.00.860.00.1060.0.
MATERIAL 3, AZ31R.
4360.00.460.00.660.00.860.00.1060.0.
MATERIAL 3, AZ31R.
4360.00.460.00.660.00.860.00.1060.0.
51250.0.1360.0.1460.0.1680.0.1680.0.
MATERIAL 4, TI-8MN.
4360.00.460.00.660.00.860.00.1060.0.
51250.0.1460.00.660.00.860.00.1060.0.
51250.0.1460.00.1662.0.1662.0.1662.0.
MATERIAL 4, TI-8MN.
5360.00.660.00.960.00.1160.0.1260.0.
71350.01860.0.2460.0.3495.0.3495.0/
THERMAL CONDUCTIVITY. RIUZIN-SEC-DEG R.
C
                                DATA NCT/10.10.3.2/
                         MATERIAL 1, 2024-T3.

10.002411.0.001283.0.001875.0.002267.0.002421.0.002450.

10.002411.0.002323.0.002189.0.002020.0.001485.

MATERIAL 2, 7075-T6.
20.001312.0.001890.0.002089.0.002284.0.002377.

30.002359.0.002240.0.002091.0.002000.0.001895.

MATERIAL 3, AZ313.

40.000901.0.000990.0.001161.0.001325.0.001470.

50.001584.0.001659.0.001680.0.001700.0.001700.
                         MATERIAL 1.2024-T3.

DATA TCONO/360.C,660.0.860.0.360.0.1060.0.

1150.U.1260.U.1360.0.1460.0.1680.0.

MATERIAL 2.7075-T6.

2360.00.660.00.760.00.860.00.960.00.

3106U.0.1260.0.1460.0.1560.0.1680.0.

MATERIAL 3.AZ313.

436U.00.460.00.650.00.860.00.1060.0.

51260.0.1460.0.1560.0.1662.0.1662.0.
3
                                EMISSIVITY.
                                DATA EMI/40*0.5/
```

```
HEAT OF FUSION.
          DATA HOF/170.0,170.0,158.0,188.0/
          DENSITY.
0
          DATA PHOM/C. 190.0.191.0.064. C. 171/
          MELT TEMPERATURES FOR MATERIAL GIVEN, DEGREES R.
        DATA TMELT/1680.0,1680.0,1662.0,3495.0/

DATA KLD/4*1/, MMAT7/4/

DATA NWL/14/

DATA WAVEB/0.2, 0.375, 0.625, 0.875, 1.125, 1.375, 1.625, 1.875,

1.225, 2.75, 3.25, 3.75, 4.25, 4.75, 5.25/

BATA PI/3.1415926535598/

END
                                                                                                                  BLOCKO
```

```
SUBPOUTING CONVEC (IFIRST, PO, TAM, VEL, TEMP, EL, XCON, HOONO, IE, 1 To, OCOM)

THIS DOUTINE CALCULATES THE CONVECTION LOSSES TO TO AIR-STREAM COOLING.

DIMEMSION TEMP(75), EL(75), XCON(75)

IF (IFIRST, GT.0) GO TO 40

TAM = TAM + 4.50.69

VC = AMAX1(VEL, 3C.)

AT = SQPT(2402.4*TS)

PAM = PO*(1.0-0.2*(VC/AT)**2)

PAM = PO*(1.0-0.2*(VC/AT)**2)**3.5

TS = TFS + .38*(TS - TFS)

XSFF = 1.15

NOTE = 1.15

NOTE = 1.15

OALQULATE CONVECTIVE HEAT FLOW.

40 HCONO.GT.0.0) GO TO 60

TRRIM = .5*(TEMP(1E) + TFS) + .22*(TR-TES)

OUT = 3.756+PAM*(TPOINTH(193.6))/TPPIM**2.5

1 (10-8.83E-6*TPOINT*).5*EL(1E)

PCON = HCONO.GT.6.0

HCONO.GSFAK*RE**0.5/XEFF

51 OCON = HCON*(TR-TEMP(IE))*EL(IE)

END
```

CONVEC

```
DIEMP CALCULATES HEAT FLOW BY CONDUCTION AND INTERNAL RAPIATION AND UPDATES TERREPORTUSES FOR EACH ELEMENT.
                   COMMON A (75), ALF (10.10), 3(122), CKD(75), CKDM(9), 3KL(75), 3KLM(0), 1 30MDAG (122), 60MD(16,9), 60MD(1622), 60MD(1622), 66(75), 5FMAT(9), 2 3LTIN, 7TIM(75), DTIME, 0TMIN, EL(75), FLUXT(15), HCONO, ID(20), 3 IEL(2), IELAY(6,3), IEMD, IMAT(75), TMELT(75), INDUT, IPLOT, ISEG(75), 4 ISTOP, JEL(3), 4CH, KCONDI(122), KCONDJ(122), KCPT(122), K32J(122), 6 KAT(3.5), KSJR, LAYM(75), MGOR, NGONDO, NEL, NELL(3,5), NELS(3), 6 KSJR, NELY(3), NAT, NOPT, NPPINT, NSEG, NTSPEC, NHLA, NHLI, PO, 7 PRINT, 0(75), JAB(3), QCOND(75), JCONV(3), QDMAX, OFLU, QR(50.3), 8 DPAD, JPEP(2.3), JRR(14), SEGL, T(75), TAM, TCUT, TEF(7), TEMP(75), COMMON, IFLUX(15), TEX, ITME, TLAY(3,5), IP(75), TR, TSPEC(3), ISTOP, 1 WEL, WAVE(50.3), WAVEL(14), WL(10.10), WT(75), X(75), X(70N(75), 1 KSPEG(4), XTO, Y(75), YSEG(4), CONDL(10.9), GPM(10.9), FMT(10.9), TSEG(4), XTO, Y(75), NCPT(9), NCT(9), NMATT, NWL, PI, PHOM(9), 1 TSEG(10.9), TCONO(10.9), TMELT(3), WAVEB(15)
                               UPDATES CRITICAL TEMPERATURE PATIO IF KC. TE. O
SET UP MATERIAL CONDUCTIVITIES AND SPECIFIC HEATS
 JU=NEL

OD 70 IE=1, IU

IF (1MELT( TE).EQ.1) GO TO 30

NM=1MAT( IE)

TEL=TEMP( IE)

IF (TEL.GE.TMELT(NM)) GO TO 27

7 CALL INT17 (7,TEL,NCT(NM),TCOND(1,NM),GONOL(1,NM),CKL(IE))

IF (KLD(NM).EQ.2) GO TO 10

CKD(TE)=CKL(IE)

GO TO 20

10 CALL INT17 (8,TEL.NCT(NM),TCOND(1,NM),CONDD(1,NM),GKD(IE))

22 IF (TEL.LT.TMELT(NM)-10.0) GO TO 25

23 CB(IE)=8.1*HOF(NM)

GO TO 30

25 CALL INT17 (11,TEL.NCPT(NM),TCP(1,NM),CPM(1,NM),CP(IE))
 GO TO 30
CALL INTIZ (11.TEL,NCPT(NM),TGP(1.NM),CPM(1.NM),GP(TE))
GO TO 30
Z7 I=NGT(NM)
GKL(IE)=CONDL(I,NM)
IF (KLD(NM).EC.2) GO TO 29
GKD(IE)=GKL(IE)
GO TO 29
Z8 GKD(TE)=CONDD(I,NM)
Z9 I=NCPT(NM)
GO TO Z8
Z0 CONTINUS
                               CALBULATE CONDUSTIVE HEAT FLOW
ICCLUDENT CONDUCTIVE HEAT FEOM

CKI = CKL(I)

CCLUDENT CONDUCTIVE CONDUCTIVE

CCLUDENT CONDUCTIVE CONDUCTIVE CONDUCTIVE

CCLUDENT CONDUCTIVE 
                                                                                                                                             1)+1MELT( J).GT.3) GO TO 83
                                                                                                                                                                                                                                                                                                                                 ICCINCKN+CONDACT
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                   ICCI
     HO CONTINUE
```

SUBSCUTINE DIENSIKO

TEMP

```
SUBSOUTINE FLUX (IFIRST.NOPI, NFLUX, TFLUX, FLUXT, QMAX, TCJT, T, GPAD)
Č
                THIS ROUTINE CALCULATES THE INSTANTANEOUS FLUX AS A FUNCTION OF TIME. THERE ARE THREE OPTIONS.
COCOCOCOCOCOCO
               IFIST - 0, FIRST PASS

1. ALL OTHER PASSES.

NOST - 1, 6000 WATT TUNGSTEN LAMP CHARACTERISTICS.

2. STEP FUNCTION.

3. ARRITARRY, POINT BY POINT (TGUT NOT USED).

NELLY, TELUX, ELUXT - DEFINE FLUX TARLE VS. TIME.

1011 - TIME AT WHICH THERMAL EXPOSURE IS CUT OFF.

T- TIME OF WHICH THERMAL EXPOSURE IS CUT OFF.
                GRA? - INSTANTANEOUS VALUE OF FLUX.
             IF (IFIRST.GT.0) GO TO 1100
IF (NOPT-2) 100,600,700
                OPTION 1.
      100 DO 200 I=2,NRISE
     130 DO 200 I=2,NRISE

II = I

IF (TRISE(I).GT.TOUT) GO TO 300

230 CONTINUE

300 CALL INTIZ(2,TOUT, NRISE,TPISE, QDRISE, QDM)

ODFALL(1) = QDM

II = II -1

DO 400 I=1,II

TFLJX(I) = TRISE(I)

400 FLUXT(I) = ODRISE(I)*ODPAT

III = II

DO 500 I=1,NFALL

III = III +1

TFLJY(III) = TFALL(I) + TCUT

NFLJY = II + NFALL

GO TO 1000
                OPTION 2.
     600 TFLJX(1) = 0.

TFLJX(2) = TCUT

TFLJX(3) = TCUT + 1.E-5

TFLJX(4) = 100.

FLJXT(1) = QDMAX

FLJXT(2) = QDMAX

FLJXT(3) = 0.
                FLJX + (4) = WFL JX = 4
                OPTION 3.
     700 004 = 0.

00 500 I=1.NFLUX

IF (FLUXT(I) .ST. 20M) QOM = FLUXT(I)

800 CONTINUE

0074T = QOMAX/QOM

00 200 I=1.NFLUX

900 FLUXT(I) = FLUXT(I) *20RAT
1000 KETJEN
```

```
1100 DO 1200 J=JL.NFLUX
IF (TFLUX(J).ST.T) 60 TO 1300
1200 GONITMUE
JU = NFLUX
0500 = FLUXT(JL)
60 TO 1400
1300 JL = J
01 = FLUX(J-1)
1100 DO 1200 J= DO 1200
1300 JL = J
1100 DO 1200 J= DO 1400
1300 JL = J
1100 DO 1200 J= DO 1400
1300 JL = J
1100 DO 1200 J=DO 1400
1300 JL = J
1100 DO 1200 J=DO 1300
1300 JL = J
1400 J= FLUX(J-1)
1400 J= DO 1400
1500 J= DO 1400
1500 J= DO 1400
1500 J=DL NFLUX (J-1)
1600 J=DL NFLUX (J-1)
1700 J=DO 1400
1700 J=DO 1
```

FLUX

```
SUBSOUTINE INTIZ (ICODE, Y, NX, XT, YT, R)

SUBSOUTINE TOTAL IS A LINEAR INTERPOLATION ROUTINE.

GIVEN A VALUE X, INTI RETURNS THE CORRESPONDING VALUE Y.

ICODE - CODE FROM CALLING PROGRAM FOR TRACEBACK IN CASE OF EFROR.

NY = THANSION OF X-TALLE AND Y-TABLE IN CALLING PROGRAM.

YT = TABLE OF X-VALUES IN CALLING PROGRAM.

R = RETURN VALUE Y.

OITHNOTON XT(1), YT(1)

IF (Y-XT(1)) 1.8.2

IF (X-XT(1)) 1.8.2

IF (X-XT(1)) 1.8.3

SUBITE(6,11) X, (XT(1), I=1,NX)

GO TO 10

4 TO 5 I=2.NX

IF (Y-XT(1)) 7.5.5

DONTINUE

PETJON

RETURN

PETJON

RETURN

PETJON

P
```

TNT1Z

```
SUBPOUTINE RPLOT (KP)
THIS POUTINE RETS UP THE PLOT FILES FOR TEMP VS TIME.
COMMON 4(75), ALF(16,16).9(122).CKD(75), CKDM(9).JKL(75).JKLM(9).
DELTIM.DITM(75).DITME.DITMIN.EL(75).FLUXT(15).HCONO.ID(20).
LILIM.DITM(75).DITME.DITMIN.EL(75).FLUXT(15).HCONO.ID(20).
LILIM.DITM(75).DITME.DITMIN.EL(75).FLUXT(15).HCONO.ID(20).
LILIM.DITM(75).MCONDIMIN.EL(75).INOUT.IPLOT.ISEG(75).
LILIM.DITM(75).KCONDI(122).KCONDJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCPJ(122).KCP
                                  IF (KP-1) 100,400,800
                            NPLOT = 0
N = (ISTOP + OELTIM)/DELTIM
NUPLOT = 2*N/NPOINT
IF (NUPLOT.EO.0) NUPLOT = 1
SET UP PLOT FOR FACH SURFACE AND LAYER BOUNDARY.
00 300 IS=1.NSE3
ILU = NLAY(IS)
IFLAY(1.IS) = IEL(IS)
00 200 IL=1.ILU
IFLAY(IL+1.IS) = IELAY(IL.IS) + NELL(IS.IL)
IELAY(ILU+1.IS) = IELAY(ILU+1.IS) - 1
IS = 1
                                 460
                                                        TA (NELL (IS, IL-1) . LE. 1) GO TO 520 = TA + .5*(TA - TEMP(IE-2)) TEMP(IE) (TA.EQ. 0.0) GO TO 580 = TB
                                                       (NELL(IS,IL).LE.1) GO TO 560

= TR + .5*(T3 - TEMP(IE+1))

(TA*TR.EQ.0.3) GO TO 500

= 0.5*(TA + TB)

(T2.GT.T1) GO TO 565
```

```
TE = ANAY1(TE,T2)
TE = AMIN1(TE,T1)
TE = AMIN1(TE,T1)
TE = AMIN1(TE,T1)
TE = AMIN1(TE,T2)
TE = AMIN1(T
```

RPLOT

```
[NE RPPINT

COMMON A(75), AL=(10,10), B(122), CKD(75), CKDM(9), CKL(75), CKLM(2),

COMMON A(75), AL=(10,10), B(122), COND)(122), CP(75), CPMAT(2),

COMMON A(122), CONDD(10,9), CONDJ(122), CPC(75), CPMAT(2),

COMMON A(122), CONDD(10,9), FLUXT(15), HCONO, IO(20),

ISL(3), IELAY(5,3), IEND, IMAT(75), IMELT(75), INOUT, IPLOT, ISEG(75),

ISTOP, JEL(3), KCH, KCONDI(122), KCONDJ(122), KCPJJ(122), KCPJJ(122),

KMAT(3,5), KSJR, LAYN(75), MCOP, NCONDC, NEL, NELL(3,5), NELS(3),

KMAT(3,5), KSJR, LAYN(75), NCONDC, NEL, NELL(3,5), NELS(3),

KMAT(3,5), KSJR, NCONDC, NEL, NELL(3,5), NELS(3,5),

KMAT(3,5), NELS(3,5), NELS(3,5),

KMAT(3,5), NCONDC, NELL(3,5), NCONDC, NELL(3,5), NE
                                              HOLIF (6,1000) TIME, GRAD, GFLU

WITT (5,1100)

OO 30 IE=1.NFL

IS = ISEG(IE)

IE1 = IEL(IS)

IF (IE.E0.IE1) 30 TO 50

IF (IE.E0.IE2) GO TO 60

WRITE (6,1200) IS.IE.TEMP(IE).QCOND(IE).Q(IE)

GO TO 90

50 IF (IE.E0.IE2) 3RER(1,IS) = ORER(1,IS) + RER(2,IS)

WRITE (6,1300) IS.IF.TEMP(IE).QAR(IS).QRER(1.IS).

1 200NV(IS).QCOND(IE).Q(IE)

60 TO 90

63 WRITE (6.1400) IS.IE.TEMP(IE).ORER(2,IS).2COND(IE).Q(IE)

CONTINUE

BETJON
        0
1000 FORMAT (13H0TIME, SFC = E14.6/

1 34H INCIDENT FLUX, BTU/IN**2/SEC = E14.6/

2 34H INCIDENT FLUENCE, BTU/IN**2 = E14.6)

1100 FORMAT (31H0 SESMENT ELEMENT TEMPERATURE, 9X,

1 37HH E A T E LOW, B T U / S E C/

2 29H NUMBER NUMBER OEGREES R, 4X,

3 49HARSORPTION PERADIATION CONVECTION CONDUCTION, 6

1200 FORMAT (1H, 16.6X, 12.4X, E11.4, 41X, E11.4, 2X, E11.4)

1300 FORMAT (1H, 16.6X, 12.4X, E11.4, 5(2X, E11.4))

1400 FORMAT (1H, 16.5X, 12.4X, E11.4, 15X, E11.4, 13X, 2(2X, F11.4))

FNO
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                               CONDUCTION 6X . 3 HMFT)
```

SUSINA

```
C
          1 FORMAT (6|12)

2 FORMAT (6|12)

3 FORMAT (6|12)

3 FORMAT (2044)

ISTOP=0

READ (5,3) (ID(I),I=1.20)

IF(ID(I).NE.IEN1.OR.ID(2).NE.IEN2) GO TO 4
    IST7P=1
GO TO 899

4 REAT(5,1) INOUT, TPLOT
IF (INOUT, GT. 0) WRITE (6.951) (ID(I), I=1,20), IPLOT
OST FORMAT (IH1,2014//30H PLOT CODE (0 FOR NO PLOTS) = I2)

MAYEL (I) = .5* (WAVER(I) + WAVER(I+1))

XMIN=10.0E10
XMIX=-10.0E10
IK=1
               TK=1
NFL=0
NMAT=NMATT
      NMAT=NMATT

KSU2=1

NS = 1

READ(5,1) NSEG

ISU=NSEG+1

DO 13 IS=1,ISU

READ(5,2) XSEG(IS), YSEG(IS)

XMIN=AMIN!(XMIN.YSEG(IS))

13 XMAX=AMAX1(XMAX, XSEG(IS))

ISJ = ISU - 1

DO 26 IS=1,ISU

READ(5,1) NLAY(IS)

ILU=NLAY(IS)

DO 14 IL=1,ILU

READ(5,1) KMAT(IS,IL), NELL(IS,IL)

READ(5,1) KMAT(IS,IL), NELL(IS,IL)

READ(5,1) KMAT(IS,IL), NELL(IS,IL)

CONTINUE
               CALCULATE SEGMENT LENGTHS
3
               SET UP ELEMENT NUMBERING SYSTEM, CALCULATE ELEMENT THICKNESS,
3
```

```
AREA, WEIGHT, LENGTH, POSITION

THICK = TLAT (LS, IL) / FLOAT (NFLL(IS, IL))

APEA = THICK + SEGL

IF (IL. GT. 1) GD TO 22

IFI = ISL(IS)

Y(IST) = 0.5* (YEEG (IS) + YSEG (IS+1) + THICK + DYL)

Y(IST) = 0.5* (YEEG (IS) + YSEG (IS+1) + THICK + DYL)

Y(IST) = 0.5* (YEEG (IS) + YSEG (IS+1) + THICK + DYL)

Y(IST) = 0.5* (YEEG (IS) + YSEG (IS+1) + THICK + DYL)

Y(IST) = 0.5* (YEEG (IS) + YSEG (IS+1) + THICK + DYL)

Y(IST) = 0.5* (YEEG (IS) + YSEG (IS+1) + THICK + DYL)

Y(IST) = 1.5* (YEEG (IS) + YSEG (IS+1) + THICK + DYL)

Y(IST) = 1.5* (YEEG (IS) + YSEG (IS+1) + THICK + DYL)

Y(IST) = 1.5* (YEEG (IS) + THICK + DYL)

Y(IST) = 1.5* (YEEG (IS) + THICK + DYL)

Y(IST) = 1.5* (YEEG (IS) + THICK + DYL)

Y(IST) = 1.5* (YEEG (IS) + THICK + DYL)

Y(IST) = 1.5* (YEEG (IS) + THICK + DYL)

Y(IST) = 1.5* (YEEG (IS) + THICK + DYL)

Y(IST) = 1.5* (YEEG (IS) + THICK + DYL)

Y(IST) = 1.5* (YEEG (IS) + THICK + DYL)

Y(IST) = 1.5* (YEEG (IS) + THICK + DYL)

Y(IST) = 1.5* (YEEG (IS) + THICK + DYL)

Y(IST) = 1.5* (YEEG (IS) + THICK + DYL)

Y(IST) = 1.5* (YEEG (IS) + THICK + DYL)

Y(IST) = 1.5* (YEEG (IS) + THICK + DYL)

Y(IST) = 1.5* (YEEG (IS) + THICK + DYL)

Y(IST) = 1.5* (YEEG (IS) + THICK + DYL)

Y(IST) = 1.5* (YEEG (IS) + THICK + DYL)

Y(IST) = 1.5* (YEEG (IS) + THICK + DYL)

Y(IST) = 1.5* (YEEG (IS) + THICK + DYL)

Y(IST) = 1.5* (YEEG (IS) + THICK + DYL)

Y(IST) = 1.5* (YEEG (IS) + THICK + DYL)

Y(IST) = 1.5* (YEEG (IS) + THICK + DYL)

Y(IST) = 1.5* (YEEG (IS) + THICK + DYL)

Y(IST) = 1.5* (YEEG (IS) + THICK + DYL)

Y(IST) = 1.5* (YEEG (IS) + THICK + DYL)

Y(IST) = 1.5* (YEEG (IS) + THICK + DYL)

Y(IST) = 1.5* (YEEG (IS) + THICK + DYL)

Y(IST) = 1.5* (YEEG (IS) + THICK + DYL)

Y(IST) = 1.5* (YEEG (IS) + THICK + DYL)

Y(IST) = 1.5* (YEEG (IS) + THICK + DYL)

Y(IST) = 1.5* (YEEG (IS) + THICK + DYL)

Y(IST) = 1.5* (YEEG (IS) + THICK + DYL)

Y(IST) = 1.5* (YEEG (IS) + THICK + DYL)

Y(IST) = 1.5* (YEEG (IS) + THICK + DYL)

Y(IST) = 1.5* (YEEG (IS) + THICK + DYL)

Y(IST) = 1.5* (YEEG (IS) + THICK + DYL)

Y(IST) = 1.5* (YEEG (IS) + THICK
                                                                                            AREA, WEIGHT, LENGTH, POSITION
3
                                                                                            CALCULATE HEAT CONDUCTION CONSTANTS FOR ELEMENTS WITHIN BRANCH
                                   CALCULATE HEAT CONDUCTION CONSTANTS FOR ELEMENTS WITHIN CONSTANTS FOR ELEMENTS 
                                                                                            ELEMENTS OF ADJACENT SEGMENTS
                                               35 RIJ=AMIN1(DBI, J3J) - AMAX1(DTI, DTJ)

KCONDJ(IK) = IEJ

KCONDJ(IK) = IEJ

R(IK) = RIJ

CONDJ(IK) = EL(IEI)/(Z. 0*BIJ)
```

```
CONDUCIK) = EL(IEJ) / (2.8*8IJ)

CONDAG(IK) = 0.0

KCPI(IK) = 1

KCPI(IK) = 1

IK=IK+1

IF (DPJ.GI.DRI) GO TO 38

IF (IFJ.NE.IELUJ) GC TO 36
            IF (IFJ.NE.IELOJ) GO 10

ISC 0

GO 10 38

36 IEJ=IEJ+1

OTJ=09J

DRJ=DTJ+T(IEJ)

IF (DTJ.LT.OBI) GO TO 35
                                            ELEMENTS OF SAME SEGMENT
          35 IF (IFI.EO.IELUI) GO TO 40

KCONDI(IK)=IEI

KCOND.(IK)=IEI+1

3IJ=EL(IFI)

3(I<)=BIJ

COND.(IK)=T(IFI)/(2.0*BIJ)

COND.(IK)=T(IFI)/(2.0*BIJ)

COND.(IK)=0.0

KCPI(IK)=2

KCPI(IK)=2

KCPI(IK)=2

IX=IK+1

73 CONTINUE

40 CONTINUE

NCOND.C = IK-1
                                               MCONDC = IK-1
                                            SET UP MATERIAL PROPERTIES TABLES FOR ADDED MATERIALS
                                     NML=NMATT+1
IF (NML.GT.NMAT) GO TO 260
70 250 NM=NML, VMAT
READ (5,1) MCOR
IF (MCOR.FD.0) GO TO 115
TF (TNOUT.EQ.0) GO TO 113
IF(MCOR.GT.NMATT) GO TO 112
WPITE (6.935) NM, PRCOP(MCOR), PRCOR(MCOR+4)
GO TO 113
WPITE (6.9351) NM, MCOR
READ (5,1) KCH
IF (KCH.FQ.1) GD TO 115
NAT(NM)=NAT(MCOR)
NTU=NAT(NM)
TF (KCH.F0.1) 30 TO 115

NAT(MM) = NAT(MCOR)

NTJ = NAT(MM)

DO 114 NT=1.NTU

TAL= (NT.NM) = EMI(NT.MCOR)

EMI(NT.NM) = EMI(NT.MCOR)

DO 114 NW=1.NWL

114 ALFAT(NT.NW.NM) = ALFAT(NT.NW.MCOR)

GO TO 120

115 DEAD (5.1) NAT(NM)

IF (INDUT.EQ.1) HPITE (6.936) PRKCH(1).PRKCH(12)

IF (INDUT.EQ.1) HRITE (6.918) NAT(NM)

DO 11A NT=1.NTJ

READ (5.2) TALF(NT.NM).EMI(NT.NM)

READ (5.2) TALF(NT.NM).EMI(NT.NM)

READ (5.2) HL(NA.MT).ALF(NW.NT)

OO 17 NW=1.NWL

CALL INTIT (12.WAVEL(NW).NWLA.WL(1.NT).ALF(1.NT).ALFAT(NT.NW.NM))

17 GOVITNUE

IF (INDUT.EQ.0) GO TO 118

WPITE (6.913) TALF(NT.NM).EMI(NT.NM).NWLA.(HL(NA.NT).ALF(NW.NT).NW

1=1.NWLA)

1=1.NWLA)
```

```
KL)(NM)=KLD(MGOR)
NTJ=NCT(NM)
DO 122 NT=1,NTJ
TCOND(NT,NM)=TCOND(NT,MGOP)
SONTL(NT,NM)=CONDL(NT,MCOP)
LF (KLO(NM).EG.1) GO TO 122
CONDD(NT,NM)=CONDO(NT,MCOP)
CONTINUE
CO 130
CONSTRUCT, NAM = CONDICAT, ACCOUNT OF THE (KICHAN), ECLIS ON 10 122 CONDICAT, NAM = CONDICAT, ACCOUNT OF THE CONDICAT, NAM = CONDICAT, ACCOUNT OF THE CONDICAT, NAM = CONDICAT, ACCOUNT OF THE CONDICAT, NAM = CONDICAT, NAM =
```

```
DETERMINE EMISSIVITIES OF ADDED MATERIALS IF REQUIRED
  220 IF (FMI(1,NM).GT.0.0) GO TO 250
NTU=NAT(NM)
DO 240 NT=1,NTU
NP=2
                 EMISS=0.0
HL1=259000.0/TALF(NT.NM)
DO 230 N=3.130
 EN=N

7ET4=0.1*FN

NL2=WL1/FN

IF (WL2.LT.WL(1.NT).OR.WL2.GT.WL(NWL4.NT)) GO TO 228

CALL INT1Z(4.WL2.NWL4.WL(1.NT).ALF(1.NT).AL)

EMISS=EMISS+AL*ZETA**3/(EYP(7ET4)-1.0)

GO IO 230

229 IF (NF.GT.C) GO TO 230
            2 PANGE PROVIDED, /44H ARSON

230 CONTINUE

EMI(NT,NM)=1.5*EMISS/PI**4

240 CONTINUE

250 CONTINUE

260 IEU=NEL

DO 270 IE=1,IEU

NM=1MAT(IE)

270 WT(IE)=4 (IE)*RHOM(NM)

DO 390 IE=1,IEJ

390 IMELT(IE) = 0
                                                                                                                                                                                                                             UNAVAILABLE
SET UP DISTANCES FROM LEADING EDGE FOR CONVECTI

ISU=NSEG

DO 480 IS=1,ISJ

IE=IFL(TS)

YCON(IE) = Y(IE)

IF1 = IE + 1

IF2 = IF + NELS(IS) - 1

IF (JEU.LT.IE1) GO TO 480

DO 460 I=IE1.IEJ

460 XCON(I) = XCON(IE)

480 CONTINUE

READ (5.2) VFL.PO.TAM.HCONO

IF(INOUT.EO.G) 30 TO 500

MRITE (6.952) VEL.PO.TAM

IF(HCONO.GT.0.0) WRITE(6.953) HCONO

500 PSAD(5.1) NOPT

IF(INOUT.GT.0) WRITE(6.954) NOPT

PSAD(5.2) TOUT

IF(INOUT.GT.0) APITE(6.955) OD MAX

IF(NOUT.GT.0) ARITE(6.956) TOUT

GO TO 600

550 PSAD(5.1) NFLUX

READ(5.2) TEUX(I).FLUXT(I)

IF(INOUT.GT.0) ARITE(6.957)

DO 586 I=1.NFLUX

READ(5.2) TEUX(I).FLUXT(I)

IF(INOUT.GT.0) ARITE(6.958) I.TFLUX(I).FLUXT(I)

600 NTSPEC=1
                  SET UP DISTANCES FROM LEADING EDGE FOR CONVECTION
                 F(INOUT.GT.0) ARITE(6,958) I.TFLUX(I), FLUXT(I)
CONTINUE
NTSPEC=1
READ(5,1) NWLT, NTSPEC
IF(NWLI.E0.0) 30 TO 700
TF(IMOUT.ST.0) MPITE(6,959) NWLI, NTSPEC
TF(NTSPEC.E0.1) GO TO 650
READ(5,2) (TSPEC(I), I=1, NTSPEC)
TF(IMOUT.GT.0) ARITE(6,960) (TSPEC(I), I=1, NTSPEC)
IF(IMOUT.GT.0) ARITE(6,961)
DO 580 J=1,NYSPEC
DO 580 J=1,NYLI
```

```
2547(5.2) WAVE(J,I),00(J,T)
TE(IMOUI.ST.2) MRITE(6,352) I,J,WAVE(J,I),00(J,I)
                                             DEATIFE, 2) DELTIM, ISTOP, DELTIM, TSTOP, PRINT TE (INCUI.GI. 0) WRITE (E, 040) DELTIM, TSTOP, PRINT WRITE (E, 963)
                                             60 10 420
300 IF (IMAT(IE).FQ.IMAT(IE-1).AND.T(IE).EQ.T(IE-1)) GO TO 820
               30 16 - 11 + 1

320 VALTE (5.202) IE, IS, IL, IMAT(IE), T(IE)
343 CONTINUE
363 SETURN
321 VALUE (6.02) IE.IS.IL.TMATTIE)
323 ORT JON
323 ORT JON
325 ORT JON
326 ORT JON
327 ORT JON
328 ORT JON
328 ORT JON
329 ORT JON
329 ORT JON
320 ORT JON
320 ORT JON
321 ORT JON
322 ORT JON
323 ORT JON
323 ORT JON
324 ORT JON
325 ORT JON
325 ORT JON
326 ORT JON
327 ORT JON
327 ORT JON
328 ORT
```

```
SUBPOUTINE SPECT (IFIRST, NTSPEC, TSPEC, NHLI, WAVE, QR, NWL, WAVEL, WAVEB, T, QRR)
                                        DETERMINE SPECTRAL DISTRIBUTION OF SOURCE AS A FUNCTION OF 14 WAVELENGTHS RETHER 0.2 MICRONS AND 5.5 MICRONS.

IFIRST - 0, FIRST PASS.

1, ALL OTHER PASSES.

NTSPEC - NUMBER OF TIMES THE SPECTRAL DISTRIBUTION IS SPECIFIED.

NULL - NUMBER OF INPUT TABLE ENTRIES DEFINING DISTRIBUTION. IF 0, USE BUILT-IN DISTRIBUTION FOR 6000 WATT TUNGSTEN LAMP.

WAVE - INPUT WAVELENGTHS.

OF - INPUT INTESTIFES.

NWL - NUMBER OF STANDARD WAVELENGTHS - PROBABLY 14.

WAVER - 15 WAVELENGTHS DEFINING BAND WIDTHS OF WAVEL.

T - TIME.

ORR - RELATIVE INTENSITY FOR 14 WAVELENGTHS.
                                        DIMENSION QR(50,3), WAVE(50,3), WAVEX(15,3), QRX(15,3), QRX(15,3),
DATA 6/1.0,4.0,2.0,4.5.1.0/

IF (IFIRST.GT.0) GO TO 600

IF (NWLI.GT.0) 30 TO 200

USE 6 KW TUNGSTEN LAMP.

NT3>FC = 1

NWL! = NWLY

DO 100 J=1.NTSPEC

DO 100 J=1.NTSPEC

OO 500 J=1.NTSPEC

OO 500 J=1.NTSPEC

OO 400 I=1.NWL

X1 = WAVFR(I)

X5 = WAVFR(I)

DO 4.0 I=1.NWL

X1 = WAVFR(I)

OO 300 K=1.5

CALL INT17(1.X1,NWLI.WAVE(1.J).QR(1.J).FI

300 FF = FF + F*C(K)

QPR(I) = FF*(X5-X1)/12.

400 QTOI = QTOI + DRR(I)

500 DPRX(J,I) = QRRX(I)/OTOI

500 DPRX(J,I) = QRRX(I,I)

SPECTRAL DISTRIBUTION VARIES WITH TIME.
                                           SPECTRAL DISTRIBUTION VARIES WITH TIME.
 GOO IF (NTSPEC.EQ.1) RETURN
IF (T.GT.TSPEC(NTSPEC)) GO TO 800
OO 700 I=1, NWL
CALL INTIT(3, T, NTSPEC, TSPEC.ORRX(1, I).QRR(I))
700 CONTINUE
RETJEN
   800
                                     DO 300 T=1, NWL
ORR(T) = GPRX(NTSPEC.I)
```

SPECT

```
SUBSTITUTE TIN

READ IN INPUT DATA.

DIMENSION (A(40))

WOIFE (6,1000)

J = 1

100 PEAD (1,1) (AA(I), I=1,40)

IF (ED(1)) 400,200

200 WPITE (6,2000) J. (AA(J), I=1,40)

WOIFE (6,2000) J. (AA(J), I=1,40)

J = J + 1

GO TO 100

400 WRITE (5,3000)

RETJON

1 FORMAT (40A2)

1 000 FORMAT (1H1,//,10X,15HT P A P - M L/

1 16H0*** INPUT DATA/)

2000 FORMAT (1X,I4,2X,40A2)

END
```

TTN

```
SUBPOUTINE TSTEP
COMMON A(75), ALF(10,10), B(122), CKD(75), CKDM(9), CKL(75), CKLM(9),
COMMON A(75), ALF(10,10), B(122), CONDJ(122), CONDJ(122), CP(75), CPMAT(9),
COMMON A(75), ALF(10,10), CONDJ(122), CONDJ(122), CP(75), CPMAT(9),
COMMON A(75), COMMON A(75), CONDJ(122), CONDJ(122), CPMAT(9),
CMAT(3,5), KSJR, LAYN(75), MCOR, NCONDC, NEL, NELL(3,5), NELS(3),
CMAT(3,5), KSJR, LAYN(75), NROR, NCONDC, NEL, NELL(3,5), NELS(3),
CMAT(3,5), KSJR, LAYN(75), NROR, NCONDC, NEL, NELL(3,5), NELS(3),
CMAT(3,5), KSJR, LAYN(75), NROR, NCONDC, NEL, NELL(3,5), NELS(3),
CMAT(4), NESS, NESS
                                                                                                                               SUBPOUTINE TSTEP
GO 10 32

CKI=CKDM(NM)

NM=IMAT(J)

IF (KCPJ( ICC).EQ.2) GO TO GO

CKJ=CKLM(NM)

GO TO 42

CKJ=CKOM(NM)

OKIJ=CONDI( ICC)/CKI+CONDJ(

OC=1.0/OKIJ

OC J)=Q( J)+OC

OC J)=Q( J)+OC

OC J)=Q( J)+OC

OC J(ICC)

OC J)=Q( J)+OC

OC J(ICC)

OC J(ICC)
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                   IE) . WT (
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                             IE) +CPMAT (N4) +0.8/Q(
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                            IF))
```

TOTEP

```
(OTIM( IF), IE=1. IU)
3240TIME INTERVAL REQUIRED BY EACH ELEMENT/
```

TSTEP

```
CALCULATES EXTERNAL RADIATION AND CONVECTION HEAT FLOWS
FOR FACH ELEMENT.
KR = 0, SETUP ONLY.
1. CONVECTION AND REPADIATION ARE CONSIDERED.
2. CONVECTION, RERADIATION, AND EXTERNAL RADIATION ARE USED.
                                 COMMON A (75), ALF (10,10), B(122), CKD(75), CKDM(9), CKLM(9), COMMON A (75), ALF (10,10), B(122), CONDI (122), CONDI (122), CPMAT(9), CONDI (122), CDMD (122), CDMD (10,20), CDMD (10,2
                                  20 CALL SPECT (1.NTSPEC.TSPEC.NHLI.WAVE.QR.NHL.WAVEL.WAVEL.TIME.QRR) CALL FLUX (1.NOPT.NFLUX.TFLUX.FLUXT.QDMAX.TCUT.TIME.QRA))
                30 ISJ = NSEG

00 30 IS=1,ISU

IF = JEL(IS)

IF (IE.EO.0) GO TO 50

INNER SURFACE.

NM = IMAT(IE)

IF (TEL.LT.TMELT(NM)) GO TO 42

I = MAT(NM)

EMISS = EMT(I.NM)

EMISS = EMT(I.NM)

EMISS = EMT(I.NM), TALF(1,NM).EMI(1,NM).EMISS)

43 QREP(7,IS) = -(1.0E-14/3.0)*EL(IE)*EMISS*TEL**4

0(IE) = 0(IE) + QREP(2,IS)

50 IE = IEL(IS)

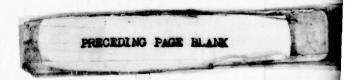
IF (IE.EG.0) GO TO 90
0
                                     CALDULATE CONVECTIVE HEAT FLOW
                                   CALL CONVEC (1,PO,TAM,VEL,TEMP,EL,XCON, HCONO,IE,TR,QCON) OCONV(IS) = QCON Q(IE) = Q(IE) + QCON
                                     CALCULATE RADIATION FROM SUPFACE ELEMENTS
                                     NM = IMAT(IE)
TEL=TEMP(IE)
IF (TEL.LT.TMELT(NM)) GO TO 62
I=NAT(NM)
EMISS=EMI(I,NM)
GO TO 63
```

SUBPOUTINE XHEAT (KE)

XHFAT

YHEAT

APPENDIX B
PROGRAM LISTING OF
APLOT



```
PROSPAM APLOT (OUTPUT, TAPE6=OUTPUT, TAPE8=514, TAPE40=514, PLOTS)
THIS OROGRAM PLOTS THE DATA GENERATED BY TRAP-ML AND STOPED ON
FILES TAPE8. THE OUTPUT PLOT TAPE IS TAPE39.
PROGRAM APLOT WAS WRITTEN BY KAMAN AVIDYNE. SEPT., 1973.
COMMON /STORE/ DAT(200,6), IDD(20), OD, TC, IL1, TIM(600)
COMMON/KACOM/NOPM, ICIN(2), ICOUT(1), INP(3), NGRAFS, NOGRAF, NCRT
DIMENSION NPC(6), XMT(2), YMT(2), XT(3), YT(4), ITZ(7), IT3(5), IPSS(6)
DIMENSION IFMT(1)
DATA NX/10/, NY/13/
DATA XT/4HTIME,4H (SE,4HC) /
DATA YT/4HTIME,4H (D,4HEG R,4H) /
DATA YT/4HTEMP,4H. (D,4HEG R,4H) /
DATA TIZ/4HPEAX,4H FLU,4HX, B,4HTU/I,4HN2/S,4HEC =,4H /
DATA TIZ/4HPEAX,4H FLU,4HX, B,4HTU/I,4HN2/S,4HEC =,4H /
DATA ITZ/4HPESAX,4H FLU,4HX,B,4HTU/I,4HN2/S,4HEC =,4H /
DATA ITS/4HPULS,4H FU,4HX,B,4HTU/I,4HN2/S,4HEC =,4H /
DATA ITS/4HPULS,3/, IC/4/, IR/5/
DATA AL/5.0/
DATA AL/5.0/
DATA NDIMEN/200/
3
                                      CALL PLOT (0.,0.,-3)
CALL FIRST (100)
3
                  50 IPLOT = IPLOT + 1

I = IPLOT

REA) (8) TIM(I), (DAT(I,J),J=1,IL1)

IF (EOF(8)) 120,76

70 WPITE (6,8) FIM(I), (DAT(I,J),J=1,IL1)

GO TO 50
          120 NC = IL1
IPLOT = IPLOT - 1
ELIMINATE MELTED REGION.
DO 200 J=1.NC
DO 150 I=1.IPLOT
N = I
3
           N = [DAT(I,J).EQ.0.0) GO TO 180
150 CONTINUE
NP3(J) = IPLOT
GO TO 200
180 NP3(J) = N-1
200 CONTINUE
SCALE THE TABLE.
                                 SCALE THE TABLE.

N=IABS(NPC(1))
CALL KALE (TIM, AL, N, 1, XM, DX, DELX)
CALL KALE (DAT, AL, N, 1, YM, DY, DELY)
IF(NC, FO, 1) GO TO 350

XMIN = YM

XMAX = XM+DX*AL
YMIN = YM

YMAX = YM+DY*AL

NLOC = 1
DO 3 CC TSC=2, NC
N=IABS(NPC(ISC))
NLOC = NLOC+NDIMEN
CALL KALE (TIM, AL, N, 1, XM, DX, DFLX)
CALL KALE (DAT(NLOC), AL, N, 1, YM, DY, DELY)
TFST = XM+DX*AL
IF(IEST.GT.XMAX) XMAX = TEST
TEST = YM+DX*AL
IF(IEST.GT.XMAX) YMAX = TEST
TEST = YM+DY*AL
IF(IEST.GT.YMAX) YMAX = TEST
CONTINUE
XMT(1) = XMIN
XMT(2) = XMAX
YMT(1) = YMIN
```

```
YMT(2) = YMAX
CALL KALE(YMT,AL,2,1,XM,DX,DELX)
CALL KALE(YMT,AL,2,1,YM,DY,DELY)
              DPAH THE AXES.
   353 XAX = 0.0

IF (YM.LT.0.0.AND.-YM/DY.LT.AL) XAX=-YM/DY

CALL KAYIS (C.C.XAX,XT,-NX,AL ,C.O,XM,DX,DELX,0)

YAX=0.0

IF (YM.LT.0.0.AND.-XM/DX.LT.AL) YAX=-XM/DX

CALL KAXIS (YAX,0.0,YT,NY,AL,90.0,YM,DY,DELY,-1)
             DPAW TITLES.

H = 0.1

CALL KASYM (0.3.5.5.H.IDD, 0.0, NT1, 3)

CALL KASYM (0.8.5.3.H.IT2, 0.0, NT2, 3)

CALL KASYM (9.9.5.3.H.OD, 0.0.1 FMT, IC, IR)

IF (NOP.ED.3) GD TO 400

CALL KASYM (0.3.5.1.H. IT3, 0.0, NT3, 3)

CALL KASYM (999.0.5.1.H. IT3, 0.0, IFMT, IC, IR)
              PLOT FACH SUPVE.
   400 NF7 = IPLOT/5
DO 500 I=1,N0
K = NDIMEN*(I-1) + 1
ND = NPC(I)
IPS = IPSS(I)
CALL KALINE (TIM(1),DAT(K),NP,XM,DX,YM,DY,NFR,IPS)
500 WPITE (6,2) I
              INDICATE MELT CONDITION.
   H = 0.1

00.600 J=1,NC

IF (NRC(J).EQ.[PLOT) GO TO 600

I = NPC(J)

K = NDIMEN+(J-1) + I

X = (TIM(I)-XM)/DX

Y = (DAT(K)-YM)/DY - .5*H

CALL KASYM (X.Y.H.IDUM.0.0.-7,3)
              CHECK FOR LAST PLOT.
    IF (NOGRAF.EQ.1) GO TO 700 NRITE (6.4) NOGRAF GO TO 800 PRITE (5.5) 370 NOGRAF=NOGRAF+1
200
              ADVANCE TO VEXT FRAME.
             CALL KAVANS (AL+3.5,6.0)
GO TO 23
              CLOSE THE PLOTTING SYSTEM.
    900 NO37AF = NOGRAF - 1
HPITF (6.6) NO37AF
CALL PLOTE (NN)
COCYC
              FORMAT STATEMENTS.
           7 FORMAT (13H10ATA PLOTTED/)
3 FORMAT (1H , 515.6, 57, 5615.6)
```

c END

APLOT

```
SURROUTINE FIRST (M)
          OPEN PLOT FILES FOR OFF-LINE PLOTS.
          M = NUMBER OF GRAPHS TO BE PLOTTED.
          COMMON/KACOM/NOPH, ICIN(2).ICOUT(1).INP(3).NGRAFS.NOGRAF.NORT OTHENSION I(2).((3) DIMENSION PLOTID(3)
          DATA I/4H( ,4H4A1)/,J/4H(A4)/,K/4H( ,4H18A4,4H)
DATA PLOTID/10H7299 BILL ,10HLEE PHONE ,10H 2721990
0
          N=IABS(M)
XM1 X = 10 + N
          YMAX = 11.4
CALL PLTID3(PLOTID, XMAX, YMAX, 1.0)
          NCRT = FLAG FOR TYPE OF PLOTTING.
REQUIRED AT SOME INSTALLATIONS FOR PROPER FILM ADVANCE
AND ORIGIN RESET.
                        0 = ORIGIN CONTROLLED BY USER.
1 = ORIGIN MOVED UP AND OVER 2.0 INCHES FOR USE WITH
QUICK PLOT ROUTINES.
          NCRT = 0
IF (M.GT.O) NCRT=1
9
          IF (M.GT.0) CALL KAVANS (3.0,1.0)
TRM 360 - AVCO.
CALCOMP MODEL 890 CATHODE RAY TUBE PLOTTER.
MAXIMUM PLOT AREA WITHOUT DISTORTION = 7.5X7.5 INCHES.
          CALL IDERME ('AVIDYNE ', 'A5', 0, T, 39)
          N=TAPS(M)
          IF M IS LESS THAN ZERO, USER IS EXPECTED TO ADVANCE FILM BEYOND AVOI ID FRAME WITH
          GALL CALCMP(0.0,0.0,0,2)
          IF JSER CALLS KAVANS ORIGIN WILL BE MOVED UP 1.05 INCHES. AND PIGHT 1.25 INCHES - INTENSITY WILL BE CHANGED BY -5.
          CHECK FOR OUTCK PLOT OR USER PLOTTING.
          MOVE PEN TO RIGHT OF 1.0. FRAME AND SET INTENSITY TO A LOWER LEVEL FOR FASTER REAM MOVEMENT.
          IF (M.GT.C) CALL KAVANS (1.05,1.25)
          SET COUNTERS.
NGSAFS = TOTAL NUMBER OF GRAPHS TO BE GENERATED.
NOSBAF = NUMBER OF MEXT GRAPH TO BE FLOTTED.
          NGRAFS=N

NGRAFS=1

NGRW=L

ICIV(1)=I(1)

ICIV(2)=I(2)

ICIV(1)=X(1)

INP(1)=X(1)

INP(2)=X(2)

INP(2)=X(3)
```

1 FOR AT (31HDPLOT ROUTINE APLOT INITIALIZED)

SET JEN

0

FIRST

```
SUISOUTINE KALIVE (KV, YV, N, XM, DY, YM, DY, K, KONES)
               CALLING AFGUMENTS -
              YV = Y COORDINATES OF DATA POINTS TO BE PLOTTED. (REAL)

YV = Y COORDINATES OF DATA POINTS TO BE PLOTTED. (REAL)

N = MUMBER OF DATA FOINTS TO BE PLOTTED. (INTEGER)

A MEGATIVE N INDICATES A POINT BY POINT PLOT WITH

NO CONNECTING LINES.

YM = MINIMUM VALUE OF VAPIABLE ON ARSCISSA. (REAL)

DX = SCALE FACIDE (UNITS/INCH) FOR ARSCISSA. (REAL)

DY = SCALE FACIDE (UNITS/INCH) OF VARIABLE TO BE PLOTTED ON

PROJUCT OF VAPIABLE TO BE PLOTTED ON OFDINATE. (REAL)

REPOSITIVE ASSECTATIVE CODE. (INTEGER)

TO BE PLOTTED. THE NUMBER OF THE CHARACTER IS GIVEN IN

KODES (1).
                                 TE NEGATIVE, THE USER SUPPLIES THE PLOT 70DE FOR HIS OWN SPECIAL CHARACTER IN KODES.

<ODES(1) = NUMBER OF WORDS DESCRIBING SYMBOL.

<ODES(2) = START OF SYMBOL.
                                 IF K = 0. (1) NO SPECIAL SYMBOL IS TO BE USED, AND

(2) EVERY POINT IN THE TABLE IS TO BE PLOTTED.

MAGNITUDE OF K INDICATES POINTS IN TABLE TO BE PLOTTED.

K = 1, EVERY POINT.

K = 3. EVERY THIRD POINT.
               DIMENSION XV(1), YV(1), KODES(1)
IF (N.EQ.O) RETURN
()()()
               CHECK FOR CONTINUOUS LINE.
               IF(V.LT. 0) GO TO 20
               7744 CONTINUOUS GURVE.
       X=(XV(1)-YM)/DX
Y=(YV(1)-YM)/DY
CALL FLOT(X,Y,3)
DO 10 I=1,N
X=(XV(I)-YM)/DX
Y=(YV(I)-YM)/DX
10 GALL FLOT(X,Y,2)
.....
                CHECK FOR SPECIAL SYNBOLS.
               TF (K.EQ.0) GO TO 50
                MONE BEN TO FIRST POINT.
       20 Y=(YV(1)-YP)/DX
Y=(YV(1)-YM)/DY
GALL PLOT(Y,Y,3)
M=[ARS(N)
                CHESK FOR USER SUPPLIED CODE.
                TF ( .LT. 6) GO TO 50
               L=<
IK=:
YOD:=-KODES(1)
       LIFT PEN
```

```
GO CALL PLOT(X,Y,3)

OF USER SUPPLIED CODE.

SO KODE=-20-KODES(1)

IK=2
L=-4
GO TO 30
END
```

KALINE

```
SUBPOUTINE KALEID, AL, NP. J. AMIN, DA, DELI
                                                 SURPOUTINE KALE PROVIDES SCALING INFORMATION FOR THE AXIS POUTINE KAXIS.
ACCEPTION OF THE PROPERTY OF T
                                               CALLING APGUMENTS -
                                                 TUPJT -
                                                                                      = ARRAY OF DATA POINTS TO BE EXAMINED. (REAL)
= AXIS LENGTH, INCHES. (PEAL)
= NUMBER OF POINTS TO BE SCANNED IN ARRAY. (INTEGER)
= JUMP CODE WHOSE MAGNITUDE DETERMINES
THE DATA TO BE SCALED.
J = 1, USE EVERY POINT.
J = 2, USE EVERY SECOND POINT.
J = 3, USE EVERY THRID POINT.
ETC.
                                                                                                         ETC.
                                                 OUTPUT -
                                               AMTY = STARTING VALUE FOR VARIABLE ON AXIS. (REAL)

DA = SCALE FACTOR, UNITS PER INCH. TO BE USED ON AXIS. (REAL)

DEL = UNITS PER TIC MARK. (REAL)

ALLOWABLE DIVISIONS - 1.0, 2.0, 5.0 * POWER OF TEN.
                                               CAUTION - WHEN PLOTTING ON GRAPH PAPER, PLEASE NOTE THAT THE NUMBER OF MAJOR TIC MARKS IS DETERMINED BY THE DATA. UNLIKE OTHER PLOT PACKAGES (SUCH AS CALCOMP, ETC.) THIS ROUTINE WILL NOT ALWAYS PRODUCE ONE MAJOR TIC MARK PER INCH.
                                                DIMENSION D(1)
3
                                                 ORTAIN TOTAL NUMBER OF POINTS IN ARRAY, N.
Ć
COO
                                               AMIN = MINIMUM VALUE OF DATA IN ARRAY.

AMIN = D(1)

AMAX = AMIN
                                                00 20 I=1.N.J
                                                 T = TEST VALUE.
                                               T=)(T)
IF (AMIN.LT.T) 30 TO 10
AMIN=T
GO TO 20
IF (AMAX.GT.T) 30 TO 20
AMAX=T
                            20 CONTINUE

IF (AMAX.EQ.AMIN) AMAX=1.001*AMIN

IF (AMAX.LT.AMIN) AMAX=0.99*AMIN

IF (AMAX.NE.AMIN) GO TO 25

IF (
                                                SPECIAL DERUG OUTPUT FOR COC 6600.
                              I=LOCF(D)
WRITE (6.2) I
2 FORMAT (28H STARTING ADDRESS OF DATA = 08/)
                                               A44×=1.0
                                                MINIMUM AND MAXIMUM SHOULD FALL ON TIC MARKS. ALLOWABLE TIC MARKS = 1, 2, 02 5. R = PANGE.
```

KALE

```
25 R=4MAX-AMIN

POL=P/AL

DIM=ALOGID(ROL)

T=INT(ROL)

TE (POL.GI.1.G) GO TO 39

DEL=10.0**I

GO 10.50

30 IF (ROL.GI.2.0) GO TO 40

DEL=2.0**IC.0**I

GO 10.50

40 IF (POL.GI.5.0) SO TO 45

DEL=5.0**IC.0**I

SO TO 50

45 DEL=6.0**(I+1)

THINT(DUM)

IF (FLOAT(I).GI.DUM) I=I-1

AMIN=FLOAT(I)*DEL

P=AMAX-AMIN

DIM = 1.001*DEL*AL + AMIN

IF (AMAX.IT.DUM) AMAX = DUM

NT = NUMBER OF TIC MARKS REYOND STARTING VALUE.

NT=INT(RYDEL-0.0001)+1

IF (NT.GI.INT(AL)) GO TO 25

DELJEN

END
```

KALE

```
SUBROUTINE KANUA (X, Y, H, F. ANGLE, IFMT, NC. NR)
              SURPOUTINE KANUM PLOTS NUMBERS ACCORDING TO FORMAT IFMT.
              CALLING ARGUMENTS -
             Y.Y = CCCRDINATES OF LOWER LEFT CORNOR OF NUMBER
TO BE PLOTTED, IN INCHES, FROM THE ORIGIN. (REAL)
H = HEISHT, IN INCHES, OF PLOTTED NUMBER.
E = NUMBER TO BE PLOTTED.
CAM BE EITHER INTEGER OR DECIMAL.

ANGLE = ANGULAR ORIENTATION, IN DEGREES, OF PLOTTED NUMBER
MEASURED COUNTERCLOCKWISE FROM HORIZONTAL AXIS. (REAL)
IFMT = FORMAT UNDER WHICH NUMBER, F, IS TO BE WRITTEN.
ANY ALLOWABLE FORTRAN FORMAT THAT CORRESPONDS
IN TYPE TO F. (HOLLEPITH)
NC = NUMBER OF CHARACTERS IN FORMAT, IEMT. (INTEGER)
NP = NUMBER OF CHARACTERS PESULTING FROM FORMAT. (E.S. E16.4)
RESULTS IN 16 CHARACTERS.) (INTEGER)
C
              COMMON/KACOM/NCPW.ICIN(2),ICOUT(1),INP(3),NGRAFS,NOGRAF.NCRT
DIMENSION NEMT(18),IEMT(11,NOUT(18)
DATA NEMT(1)/44( /.NEMT(18)/44) /
3
              IF (NO.EC.0) GO TO 20
              N = NUMBER OF ALPHAMERIC WORDS IN IFMT.
              N=NO/NOPW.LT.NO) N=N+1
              FILL FORMAT APRAY.
              00 1 T I = 1 N (T)
      13 J=J+1
IF (J.LT.18) NF4T(J)=NFMT(18)
              CONVERT TO ALPHAMERIC.
             WPITE (40, NFMT) F
REATIND 40
N=N7/NC>W
IF (N*NC>W.LT.N>)N=N+1
PF4) (40, TNP) (NOUT(I), I=1, N)
PF4IND 40
              PLOT CODE.
              CALLEN
                          KASYM (X,Y,H,NOUT,ANGLE,NP,3)
```

KANUM

NO CONTRACTOR OF THE

SUBROUTINE KASYY (X3, Y3, HEIGHT, IBCD, ANGLE, NCHAR, IPENC)

SUBROUTINE KASYM DRAWS ALPHNUMERIC INFORMATION AND SPECIAL SYMBOLS AT ANY ANGLE AND ANY SIZE ON THE PLOTTING AREA.

IN ITS STANDARD FORM SYMBOL IS USED TO PRINT TEXT MATERIAL ALONG AXES, AND TITLE GRAPHS. THE LETTERS A THROUGH 7, DIGITS O THROUGH 9 AND ALL OTHER STANDARD FORTRAN CHARACTERS ARE AVAILABLE.

IN AN ALTERNATE FORM, IT WILL PLOT A CENTERED SYMBOL FROM THE SPECIAL SYMBOL TABLE, OR A USER SUPPLIED SYMBOL.

THE ARGUMENT NOTAR CONTROLS THE FORM TO BE USED.

APSIMENT LIST
X3, Y3 = COORDINATES, IN THICHES, OF EITHER

(1) THE LOWER LEFT CORNER OF THE

FIRST CHARACTER OF STANDARD TEXT TO

RE PRINTED - 02
(2) THE CENTER OF THE SPECIAL SYMBOL

THE TEXT CHARACTERS, OP SPECIAL SYMBOLS

TO BE PLOTTED.

IF HEIGHT IS GREATER THAN 0.0, THE UNITS

ARE ASSUMED TO BE IN INCHES.

IF HEIGHT IS LESS THAN 0.0, THE UNITS

ARE ASSUMED TO BE IN CENTINETERS.

TEXT MATERIAL TO BE PRINTED, OR SPECIAL SYMBOL CODE, DEPENDING UPON THE VALUE OF NCHAR.

STANDARD ALPHANUMERIC TEXT (BCD OR A-TYPE FORMAT) SHOULD BE LEFT-JUSTIFIED AND CONTIGUOUS IN (1) A SINGLE VARIABLE, (2) AN ARRAY, OR (3) A HOLLERITH LITERAL (IF THE COMPILER PERMITS). SPECIAL SYMBOL CODES SHOULD BE INTEGER FORMAT.

= ANGULAR ORIENTATION, IN DEGREES, AT WHICH CHARACTER LINE IS TO BE PRINTED. POSITIVE VALUES ARE MEASURED COUNTERCLOCKWISE FROM THE HORIZONTAL X-AXIS, AND NEGATIVE VALUES ARE MEASURED CLOCKWISE. ANGLE

NCHAP = CONTROL CONSTANT FOR TYPE OF MATERIAL TO BE PRINTED ON PLOT.

A POSITIVE NCHAR INDICATES STANDARD ALPHANJMERIC TEXT IS TO BE PRINTED. NCHAR IS THE NUMBER OF CHARACTERS TO BE PRINTED.

A NEGATIVE NCHAR INDICATES A SPECIAL CENTERED SYMBOL IS TO BE PLOTTED. IF THE ABSOLUTE VALUE OF NCHAR IS BETWEEN 1 AND 14 A SPECIAL SYMBOL FROM THE TABLE IS PRINTED. IF THE ABSOLUTE VALUE OF NCHAR IS GREATER THAN 14. THE USER SUPPLIES HIS OWN SPECIAL SYMBOL IN IBCO. ARE REQUIRED FOR THE MORE THAN THO WORDS ARE REQUIRED FOR THE DIMENSION OF THE NEW SYMBOL, EXPAND THE DIMENSION OF THE NEW SYMBOL IN THE CALLING PROGRAM IN NOW ASSUMED TO BE TWO.

= PEN CODE. = 2, PEN DOWN DURING MOVE TO (X3, Y3). = 3, PEN UP DURING MOVE TO (X3, Y3).

```
NOTAL CODES FOR STANDARD FORTRAN CHARACTERS ARE
CONTAINED IN NORMAL, AND SPECTAL SYMBOLS IN NODES.
NSYM IS THE NUMBER OF SPECIAL SYMBOL CODES IN NODES.
        COMMON/KACOM/NOPW.ICIN(2).ICOUT(1),INP(3),NGPAFS.NOGRAF.NCRT DIMENSION ICHARS(10)  
DIMENSION I3CD(1),IN(14),INORM(51),NODES(90),NORMAL(205)  
DIMENSION NORM(51)  
DATA YSAVE/0.G/
      C
3
         0414 LIFT/40/.NEXT/44/
0414 PI0180/1.745329252E-02/
         CHECK FOR EMPTY RUN.
         IF (NCHAR.EQ.0) RETURN
        ¥1=x3
IF (¥3.E0.999.0) X1=XSAVE
         Ŷ1=Y3
IF (Ÿ3.E0.999.0) Y1=YSAVE
         MOVE PEN TO (X1.Y1).
         CALL PLOT (X1, Y1, TPENC)
         CHANGE ANGLE TO RADIANS.
         T=ANGLE*PIO180
         INITIALIZE.
         SINT = SIN(T)
```

```
H= HF I GHT
CCC
        CHANGE CENTIMETERS TO INCHES IF NECESSARY.
        IF (H.LT.0.0) H=-H+2.54
        DEL = DISTANCE IN INCHES BETWEEN NODES.
        DEL=H/7.0
DELC=DFL*COST
OFLS=DEL*SINT
TSTAPT=1
CCC
        DETERMINE TYPE OF CHARACTERS TO BE PLOTTED.
        TE (NCHAP.GT.O) GO TO 70
CCC
        SPECIAL SYMBOL.
        N=-VCHAR
X2=X1+H+COST
Y2=Y1+H+SINT
        IF (N.GT.NSYM) 30 TO 60
         SYMPOL IS CONTAINED IN TAPLE.
    INDEX=IN(N)
INDEX=NODES(INDEX)
IPART=1
CCC
        PLOT MODES.
    20 NODF=IGHAR/100
IF (IPART.E3.2) NODE=IGHAR-100*NODE
CCC
        IF NODE = 40, LIFT PEN AND GO BACK FOR ANOTHER POINT.
        IF (NODE.EQ.LIFT) GO TO 50
CCC
        IF NODE = 44. END OF THIS SYMBOL.
        IF (NODE.EQ.NEXT) GO TO 170
         SEPARATE NODE INTO (X,Y) COORDINATES.
        IX=NODE/10
IY=NODE-10*IX
IF (IX-GT-4) IX=4-IX
IF (IY-GT-4) IY=4-IX
000
         ORTAIN COORDINATES OF NODE.
        X=X1+DFLC*FLOAT(IX)-DELS*FLOAT(IY)
Y=Y1+DELC*FLOAT(IY)+DELS*FLOAT(IX)
000
         PLOT POINT.
    IF (ISTART.EQ.1) CALL PLOT(X,Y,3)
ISTART=ISTART+1
IPEN=2
30 GALL PLOT (X,Y,IPEN)
COC
        ORTAIN NEXT POINT.
        IPART=IPART+1
IF (IPART.LE.2) GO TO 20
INDEX+1
GO TO 10
IPART.LE.2) GO TO 20
IPART.LE.2
         ISTART=1
X=X1
        9= 91
50 To 30
```

```
SPECIAL SYMBOL - USER SUPPLIED.
     60 INTEX=NUODES+1
    J=N-20

00 65 I=1.J

K=NNODES+I

65 NODES(K)=IRCD(II
3
          GO TO 13
          STANDARD FORTRAN SODE.
    70 NW=1
NC=1
          SHIFT TO CENTER OF GRID.
    X2=X1+3.6*(DELC-DELS)

Y2=Y1+3.6*(DELC-DELS)

DC7=7.0*DELC

DS7=7.0*DELC

R0 LETTER=1

TD42T = 1
          OBTAIN CODE OF CHARACTER.
      WRITE (40, ICOUT) IRCO(NW)
REAT (40, TCIN) (ICHARS(I), I=1, NCPW)
REAT (40, TCIN) (ICHARS(I), I=1, NCPW)
REAT (40, TCIN) (ICHARS(I), I=1, NCPW)
REAT (ADDE = ICHARS(LETTER)
DO RETTEN = 1, NNORM
IF (NODE, EQ. NORM(ITEN)) GO TO 96
RETTEN = 1, NODE, TRCD(NW)
1 FORMAT (66HD****AARNING**** NON-STANDARD FORTRAN SYMBOL ENCOUNTERED 1 97 KASYM. A1, 2X, A10/)
RETTEN
        1 SEL TON
          ORTAIN STAPTING POINT IN TARLE.
     95 INDEX = INCRM(ITEM)
          ORTAIN CHARACTER CODE FROM TARLE.
   100 ICHAR=NORMAL(INDEX)
          PLOT NO DE.
   110 NOTE = [CHAR/100 NODE=ICHAP-100*NODE
          IF NODE = 40. LIFT PEN AND GO BACK FOR ANOTHER POINT.
          IF (NODE.EQ.LIFT) GO TO 140
          IF NODE = 44. END OF THIS CHARACTER.
          IF (MODE. FQ. NEXT) GO TO 150
          SEPARATE NODE INTO (X,Y) COORDINATES.
          IX = MODF/1G
IY = NGOF-10+IX
IF(IY.GT.4) IX = 4-IX
IF (IY.GT.4) IY = 4-IX
          PLOT NODE.
          Y=Y2+DELC*FLOAT(IX)-DELS*FLOAT(IX)
Y=Y2+DELC*FLOAT(IY)+DELS*FLOAT(IX)
```

KASYM

```
0000000 0 0000 000000000 0
```

```
SUBROUTINE KAVANS(X,Y)

SUBROUTINE KAVANS ADVANCES THE FILM TO THE NEXT FRAME, AND PERSON THE LOSICAL ORIGIN PREVIOUSLY DEFINED.

A CALL WITH EITHER X OR Y = 909.0 WILL RESULT IN THE PREVIOUSLY DEFINED VALUES OF X AND Y BEING USED.

DATA XSAVE/0.0/, YSAVE/0.0/

X1=X
IF (X1.F0.999.0) X1=XSAVE
XSAVE=X1
IF (Y1.E0.999.6) Y1=YSAVE
YSAVE=Y1

FOR CALCOMP PEN PLOTTERS, USE

CALL PLOT(X1,Y1,-3)
FOR CATHODE RAY TUBE PLOTTERS, USE

CALL CALCMP(0.G.C.0,C000.2)
CALL CALCMP(X1,Y1,-6.3)

TO ADVANCE THE FILM, AND TO PESET THE LOSICAL ORIGIN.

RETJEN
END
```

KAVANS

SUBPOUTINE KAXIS(X,Y,LABEL,NC,AL,ANGLE,AMIN,DA,DEL,ITUPN) SUBROUTINE KAXIS DRAWS A LABELED AXIS.

CALLING ARGUMENTS
X = X-COORDINATE IN INCHES OF THE START OF THE ORIGIN. (PEAL)

Y = Y-COORDINATE IN INCHES OF THE START OF THE ORIGIN. (PEAL)

LABEL = ALPHAMERIO TEXT TO BE PRINTED ALONG AXIS. (HOLLEPITH)

NC = NUMBER OF CHARACTERS IN LABEL (INCLUDING ALL BLANKS).

A POSITIVE VALUE WILL PLACE THE TITLE ON THE

COUNTERCLOCKWISE SIDE OF THE AXIS. USUAL Y-AXIS.

A NEGATIVE VALUE WILL PLACE THE LABEL ON THE GLOCKWISE

SIDE OF THE AXIS, USUAL X-AXIS.

(INTEGER)

ANGLE = ANGULAR OPTENTATION OF AXIS. DEGREES.

(INTEGER)

AMIN = MINIMUM VALUE OF VAPIABLE ON AXIS.

VALUE RETURNED BY KALE. (REAL)

DA = SCALE FACTOR, UNITS/INCH, OF VARIABLE ALONG AXIS. (PEAL)

ITJRN = NUMBER OF 90 DEGREE ROTATION OF NUMBERS INDICATING

VALUE OF MAJOR TIC MARKS. (INTEGER)

COMMON/KACOM/NCPW, ICIN(2), ICOUT(1), INP(3), NGRAFS, NOGRAF, NGRT DIMENSION LABEL(1), JEMT(12), IEMT(3)

DATA JEMT/4HF6.1, 4HF6.2, 4HF6.3, 2HI2 ,4H, 1HX, 4H2H10, 2HI3 , 12HI5, 2HI6, 2HI4, 2HI3, 2HI1/

HGIVUM = HEIGHT OF NUMBERS ALONG AXIS. INCHES.

DSTO = DISTANCE BETWEEN AXIS AND TOP OF NUMBERS, AND BETWEEN NUMBERS AND AXIS LABEL.

HGTSYM = MAXIMUM ALLOMARLE HEIGHT OF SYMBOLS IN AXIS TITLE.

INCHES.

DATA HGTNUM/0.100/, DST0/0.09/, HGTSYM/0.130/

CENTER = NUMBER CENTERING VARIABLE.

POJALEM -

SOME SYSTEMS RIGHT JUSTIFY NUMBERS WRITTEN UNDER FORMAT STAFFMENTS. OTHERS LEFT JUSTIFY THESE NUMBERS. THUS, ON MACHINES THE MAJOR TIC MARK VALUES MAY BE SHIFTED EITHER LEFT OR RIGHT OF WHERE THEY SHOULD BE.

SOLITION -

CENTER PERMITS SENTERING NUMBERS FOR PARTICULAR SYSTEM IN USE.

CENTER = 0.0, LEFT SIDE OF NUMBER WILL LINE UP WITH MAJOR TIC MARK CENTER = 1.0, RIGHT SIDE OF NUMBER WILL LINE UP WITH MAJOR TIC MARK CENTER = 0.5, MIDDLE OF NUMBER (AS FORMATTED) WILL LINE UP WITH MAJOR TIC MARK. THUS, WITH AN IS FORMAT. THE THIRD SPACE WILL LINE UP WITH THE TIC MARK.

DATA CENTER/0.5/

SIDE=1.0 IF (NC.LT.0) SIDE=-1.0 DST=DST0 IF (SIDE.EQ.-1.0.AND.ITURN.EQ.0) DST=DST+HGTNUM CIV = 0.0 IF (ITURN.NE.0) CIT = -1.0 IF (ITURN.NE.0) (21-20)
XOR=1.0
IF (SIDE.EO.FLOAT(ITURN)) XOR=0.0
THETA=ANGLE*0.0174532952
SINT=SIN(THETA)
COST=COS(THETA)
T=ANGLE+FLOAT(ITURN)*90.0

ZIXAX

```
TM=1 * 0. 01745 32952
CTN= COS (TN)
STN= SIN (TN)
            NUMBER ORIENTATION CORRECTIONS.
            CON= SIDE*XOR*SINT*CIT-CENTER*COST*(CTT+1.0)
CON1=SIDE*XOR*DOST*CIT+CENTER*SINT*(CIT+1.0)
            PLOT MAJOR TIC MARKS.

PLOT MAJOR TIC MARKS.

MT = NUMBER OF TIC MARKS.
            AMAX = AMIN+AL *DA
            ATAX = ATIN+AL * DA

R= ATAY - AMIN

NT= [NT (P/DEL+3.1)

YR= Y

YR= Y
OCC
            TIO HEIGHT = 0.1 INCHES.
            XA=X+0.1*SIDE*SINT
YA=Y-0.1*SIDE*SOST
            YC = CONSTANT INCREMENT ALONG X AXIS.
            YC=35[*COST/DA
YC=36[*SINT/DA
CALL PLOT(X4,Y4.3)
CALL PLOT(X4,Y4.3)
J=1

NT 4= NT

5 DO 1 [ I=1.NTM

CALL PLOT(X3, Y3, 2)

XD=XC+XC
      YA=YA+YA

OALL FLOT(XB,Y3,2)

XA=XA+XA

YA=YA+YA

10 CALL FLOT(XA,YA,2)
            PLOT MINOR TIC MARKS, 16/INTERVAL.
            NT = 16 * NT
IF (J.EQ. 2) GO TO 20
             J= 2
      J=2

XA=XB+0.05*SIDE*SINT

YA=Y2-0.05*SIDE*COST

YC=-0.1*YC

YC=-0.1*YC

YC=-10.0*YC

YC=-10.0*YC
            WEITE VALUE AT TIC MARKS.
            MT=NT+1

A=475(AEL)

TF (A.LT.1.0.03.4.GF.10800.0) GO TO 60
            INTEGER FORMAT.
            J=[4] (AMIN)
J=[4] (DEL)
      IC=2

IF=1 (1)=JFMT(8)

IF (AMIN.NE.-10060.0) GO TO 24

IP=5

IFM((1)=JFMT(9)

30 TO 25

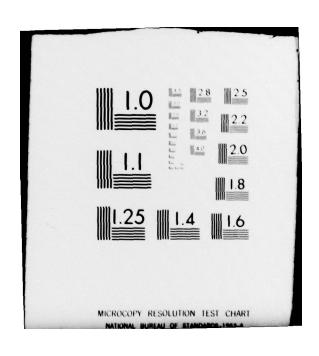
24 R=138(AMAX)
```

AD-A074 887

KAMAN AVIDYNE BURLINGTON MA
TRAP-ML-A TWO DIMENSIONAL THERMAL RESPONSE CODE TAILORED FOR TH--ETC(U)
NOV 78 W N LEE
KA-TR-154

DNA-4770F

END
PARE
PRIME
PRI



```
IF (R.LT.ABS(AMIN)) B=ABS(AMIN)
          3=8+0.1
IF (B.GE.1000.) GO TO 25
         IF (8.6E.1000.) GO TO 25
IFMI(1)=JFMT(10)
IF (8.6E.100.0) GO TO 25
IP=3
          ÎFMÎ (1) = JFMÎ (11)
ÎF (AMÎN.LÎ.0.0 .OR. DEL.LÎ.0.0) GO TO 25
ÎP=2
          TENT (1) = JEMT (4)

IF (P.GE. 10.0) 30 TO 25

IFMT (1) = JEMT (12)
 LABEL AXIS.
         S=FLOAT(NC)+SIDE+1.0
         HEALIS
IF (H.GT.HGTSYM) HEHGTSYM
 TD=4*S

XB=X+0.5*(AL-TD)*COST-SIDE*SINT*(HGTNUM+2.0*DST0+0.5*(1.0-SIDE)*H)

IF (ITU?N.NE.0) XB=XB-(FIR+DST0)*CTN

YB=Y+0.5*(AL-TD)*SINT+SIDE*COST*(HGTNUM+2.0*0ST0+0.5*(1.0-SIDE)*H)

IF (ITURN.NE.0) YB=YB-(FIP+DST0)*STN

I=IARS(NC)

CALL KASYM(XB,Y3,H,LABEL,ANGLE,I,3)

45 PETJPN
53 B=4 OG10(A)

IF (P.GT.0.0) 3=P+0.02

I=INT (B-0.01)

IF (I.GT.C.OR.I.LT.-2) GO TO 80

IFMI(1)=JFMI(3)

IF (P.LE.-1.0) IFMI(1)=JFMI(2)

IC=+

IP=5

A=AMIN

FIR=6.0*HGTNUM

X3=Y-SIDE*DST*SINT+FIR*CON

Y9=Y+SIDE*DST*COST-FIR*CON1

70 DO 72 I=1,NI

CALL KANUM(XB,YB,HGTNUM,A,T,IFMI,IC,IR)

X3=X9+XC

Y9=YB+YC

72 A=A+DFL

GO 70 40
         DESIMAL NOTATION.
         EXPONENTIAL FORMAT.
 80 A=10.0**I

AM=4 MIN/A

ADA=DEL/A

NT=VT-1

IF (ADA.GE.0.01) GO TO 90
        IF (ADA.GE.0.01) GO TO 90

I=I-1

GO TO 80

IF4Y(1)=JFMT(2)

IF (ADA.GE.1.0) IFMY(1)=JFMT(1)

IC=4

IR=5
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